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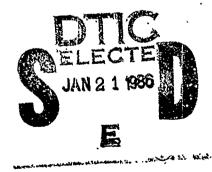
SIMULATION OF THE STATE OF THE M42/M46 GRENADE DURING PRESS LOADING

GEORGE J. SCHLENKER

DECEMBER 1985

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AMSMC/RD/MR-7

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This report is a contribution to an ongoing study of	of explosive incidents (blows)

This report is a contribution to an ongoing study of explosive incidents (blows) which occur during press loading of the M42/M46 grenade. One of the causal mechanisms posited for the blows is brittle fracture of the grenade body. The physical state of the grenade during final consolidation is of critical importance for this mechanism. A detailed simulation of the compression phase of final loading helps to verify the feasibility of this causal mechanism and to suggest means of minimizing a part of the rate of incidents. This report

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describes a continuous simulation of several, related phenomena which occur during final consolidation. These phenomena include: (a) compaction of the bulk HE, (b) production of elastic strain in the grenade body, (c) elastic compression of RDX particles in the HE, (d) kinematics of the punch, (e) heating of the HE, and (f) diffusion of heat within the grenade. A sample of simulation results is given in graphical form, with key variables displayed as functions of punch displacement and of time. Sensitivity of results to certain parameters is shown. Comparisons are made between some experimental measurements and results of the simulation. These comparisons demonstrate the validity of the simulation within the limits imposed by its scope. For the interested analyst, the implementing computer program is listed and explained.

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EXECUTIVE SUMMARY

This report is a contribution to an ongoing study of explosive incidents (blows) which occur during press loading of the M42/M46 grenade. High explosive (HE) Composition A-5 is pressed into a steel body (either M42 or M46) of the grenade in two operations: pre- and final -consolidation. The latter is nearly always where blows occur. Several causal physical mechanisms for blows have been hypotnesized. One of these mechanisms involves brittle fracture of the grenade body. If a crack opens and propagates, the rapid release of elastic strain energy at the interface of the body wall and HE fill appears to be capable of igniting the HE. The physical states of the grenade body and of the HE during the compaction phase of final consolication are critical to this mechanism. Quantification of these states helps to verify the feasibility of the brittle-fracture mechanism and to suggest means of minimizing the portion of the blow rate contributed by this mechanism. This report describes a continuous simulation of several, related physical processes which occur during the final consolidation of the HE in the M42/M46 grenade. Typical simulation results are presented in graphical form, with pertinent variables displayed as functions of both punch displacement and time. The sensitivity of these results to certain parameters is shown. For example, punch travel and work done by the punch are shown to be sensitive to the initial, preconsolidated system state. Also the max noop stress and noop strain energy are sensitive to peak punch pressure. Where possible, comparisons are made between experimental data and results of the simulation. These comparisons demonstrate the validity of the simulation within the limits imposed by its scope. For the interested analyst, the implementing computer program is listed and explained.

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AMSMC-RDA-S December 1985

MEMORANDUM REPORT

SUBJECT: Simulation of the State of the M42/M46 Grenade During Press Loading

1. Foreword

Order of topics in this report represents increasing detail. Background information is presented first to set a context and to motivate the work presented here. The goals of the subject simulation are given next, followed by the scope of— and assumptions for—the simulation. Simulation results are given in tables and graphs, followed by comments about these results. Sections on validation and methodology complete the body of the report. A source program listing of the simulation is given in Annex A. Annex B contains a memorandum concerning a model of the bulk density of Comp A-5. The following references are cited through the report.

2. References

- a. MFR, AMSMC-RDA-S, 22 Oct 85, subject: Evidence for a Steel-Supplier Effect on the Rate of Press Blows in M42 and XM77 Grenades.
- b. Technical Report ARLCD-TR-79002, Collett, R.W. and England J.T., ARRADCOM, LCWSL, Dover, NJ, Jun 79, title: Press Loading Encident Investigation of M42/M46 Grenade Bodies.
- c. MFR, AMSMC-RDA-S, 9 Aug 85, subject: Some Observations About the Explosive Sensitivity of Comp A-5.
- d. MFR, AMSMC-RDA-S, 28 Aug 85 (Revised 24 Sep 85), subject: Predicted Bulk Densities of Comp A-5 and Comp A-4 as Functions of Peak Consolidation Pressure. (Given in Annex B.)
- e. Technical Report #371, 23 Aug 85, Day & Zimmerman Kansas Div., subject: Investigation of Effects of Charge Density on Penetration in M42/M46 Grenades.
- f. MFR, DRSMC-SAS (R), 4 Nov 83, subject: Particle-Size Distribution of Nominal Class 1 RDX Before Incorporation and After Extraction From Extrusions of Comp C-4.
- g. Reference Book, Kaye, S.M., US ARDC, Dover, NJ, c. 1980, title: Encyclopedia of Explosives and Related Items, Vol 9.
- h. Text Book, Timoshenko and Goodier, McGraw-Hill, NY, c. 1951, title: Theory of Elasticity.
- i. MFR, AMSMC-RDA-S, 23 Aug 85, subject: Time Series Analysis of the Peak Consolidation Load for Pressing of HE into M42/M46 Grenades.

3. Background

Previous studies of explosive incidents during press loading of M42/M46 grenades posited various physical mechanisms for initiating an explosion. Ref 2a presents arguments supporting the mechanism of brittle fracture of the grenade body. The scenario for an explosion with this mechanism is as follows: During final consolidation of the HE (Comp A-5), a latent crack suddenly opens and propagates. A rapid release of elastic strain energy in the grenade body is deposited at the surface of the HE in proximity to the crack. This energy release over a small area is regarded as sufficient to initiate an explosion of the HE *. If this mechanism is responsible for some of the incidents, one would expect to see the observed variation in blow rate between bodies by different body producers, since variation in metal parts quality between producers is quite evident. Ref 2b also presents evidence for a steel-supplier effect, which, again, is germane to the mechanism of brittle fracture. A similar phenomenon, occurring in the powder nest, is one of the mechanisms hypothesized in Ref 2b. The very limited set of experiments reported in Ref 2b failed to demonstrate a blow with either a cracked body or a cracked nest. However, these negative results are not persuasive, because only rare and special conditions -- very rapid energy deposit in a small area -must exist in order to provoke a blow. Based on simulated results, brittle fracture must still be considered a credible mechanism.

- 4. Considering this mechanism, several actions to reduce the frequency of press blows are appropriate: (a) Improve the specified quality, including fracture toughness, of steel used in making grenade bodies. This action is suggested in Ref 2a. (b) Improve quality of grenade bodies from the "marginal" body producers. Actions started by AMCCOM Quality Assurance in 1985 address this issue. (c) Use a less sensitive HE fill, such as Comp A-4. This idea is still on the back burner. (d) Load at a lower peak consolidation pressure. This suggestion is made in Ref 2c, and justified by data which show reduced explosive sensitivity when grain breakage is minimized. Altho easy to implement, the last suggestion has met resistance on the grounds that effectiveness (penetration & lethal area) would suffer. Nevertheless, some tradeoff here seems possible to reduce blow rates.
- 5. Because of the importance of the state of stress of the grenade body to the brittle fracture mechanism, some effort was made to take experimental measurements (for example, Ref 2b). And, because of the importance of HE density to grenade performance, various data nave peen gathered relating HE density to peak consolidation load (or pressure), under both quasi-static (very slow rate of loading) and transient conditions (which occur in load plants). Some of these experimental results are compared with simple mathematical models in Ref 2d. Ref 2e presents additional results for the transient loading condition, in a study designed to relate average HE density to grenade penetration performance. Notwitnstanding these experiments to understand the processes occurring during press loading, to my knowledge no unified mathematical model has been developed to simulate the physical processes which interact during final consolidation. This memorandum presents such a model. Tho the simulation is not very sophisticated, model output is in good agreement with pertinent experimental data.

^{*} Ref 2g indicates that the threshold value of energy per unit area for shock ignition of RDX is about 10 cal/sq cm, a value which may be achieved by release of elastic strain energy in the body.

6. Goals of the Simulation

The primary goal of the simulation is to achieve better understanding of the relationship of several physical parameters pertinent to loading of the M42/M46 grenade. Understandings of this sort ultimately guide design of experiments and help to quantify improvement in rate of press blows, expected to accompany loading at lower peak pressure. A partial list of questions addressed by the simulation may clarify my intentions:

- (a) How does the HE respond to punch displacement during consolidation? For example, how does punch load vary with punch displacement and with time? (In this case, time is simply a convenient variable for use in comparison with experimental results. Since punch acceleration is quite low, inertial forces are not important. The phenomena are mostly kinematic rather than dynamic.)
- (b) How does the work done by the punch vary as a function of punch displacement and as a function of time?
- (c) What portion of the punch work is invested in elastic, i.e., recoverable, energy components? This question has a bearing on volume change in the HE following punch withdrawl. Also, these elastic strain-energy components would become sources of energy for initiating an explosion of the HE, if a metal failure occurred during compression.
- (d) Are physically significant temperature gradients developed within the RDX particles during consolidation?
- (e) Do any of the variables of interest display a large sensitivity to the initial conditions for final consolidation, i.e., does variation in preconsolidation affect final results?
- (f) How well does the model match the experimental load-versustime data?
- It is beyond the scope of this model to address issues such as:
- (a) Specifically, what happens if a brittle fracture on a grenade body occurs during consolidation?
- (b) What stress concentrations occur in various regions of the body?
- (c) Where does RDX grain fracture occur during loading, and what is its extent?
- (d) What is the nature of the flow in the HE during consolidation?

7. Scope and Assumptions

In terms of fidelity to geometric details of the grenade body, 'he model is quite crude. For the purpose of calculating hoop stress and hoop strain in the body, the grenade is modeled as a regular cylindrical sleeve, supported on one end by a smooth, unyielding surface. The HE configuration is assumed to have cylindrical symmetry

within the cavity. The simulation model is descriptive; it uses semiempirical results where appropriate, and does not insist on starting from first principles. In spite of evidence of RDX grain fracture, cited in Refs 2c and 2f, the model ignores this phenomenon. Further, because of its simple geometry, no attempt is made to describe the actual flow of the HE with respect to the punch and body cavity in an actual M42/M46 grenade. However, the effects of a pressure gradient within the HE are considered. Most of the mechanical and thermal material characteristics were obtained from standard references such as Ref 2g. The compaction model (avg HE density versus peak pressure) uses the functional form described in Annex B. This model assumes a constant pressure gradient within the HE. The pressure decreases linearly from a maximum value of pmax, at the punch face, to a minimum value of a constant, g, times pmax. The value of g used here (0.6) is obtained from data in Ref 2e. The assumed constant pressure gradient within the HE is also used in deriving an expression for the RDX elastic strain energy. Hoop stress in the simulated grenade body is calculated using the Lame equation for thick-walled cylinders (Ref 2h). A formula for body strain energy is derived from the last equation, which assumes elastic behavior.

8. Results of the Simulation

The loading simulation is implemented by a computer program COMPACT. A source listing of this program, in Simscript 2.5, is located in Annex A. The logic of this program is sketched below (p. 19, Methodology). A portion of the output from COMPACT is shown in Table 1. After echoing the input parameters, the state variables, which characterize the HE and the grenade body, are printed as functions of the displacement of the punch. The initial, average density of the HE is chosen by the program user. This value suffices to initialize the simulation for the final consolidation cycle. Punch displacement and cycle time are taken as zero at this point. Time is calculated from punch displacement, since these are kinematically related by the punch cam shape and operational speed of the press. Program variables are plotted as functions of punch travel and time. These are shown in Figures 1 thru 19, on pages 7 thru 16. Results are discussed in paragraph 10, page 17.

9. Parametric Analyses

Two simulated body dimensions were treated as parameters for sensitivity analysis: (a) the thickness of the wall of the sleeve, representing the grenade body, and (b) the effective length of the sleeve which is exposed to internal pressure. The last parameter is determined by shape of the punch and by the 1-D nature of body stress calculations. Three values of wall thickness -- 0.110 inch, 0.115 inch, and 0.120 inch -- are used at a effective body length of 1.4 inch. And, two values of the body length -- 1.0 inch and 1.4 inch -- are used with a nominal wall thickness of 0.115 inch. Results are shown in Tables 2 and 3. In each case the Initial punch pressure is 5.1 kpsi and the final pressure is 25 kpsi (to terminate the comression phase). The initial pressure corresponds to an average HE density of 1.44 g/cc. Since the final state is prescribed, certain variables, such as punch work and loading time, are not very sensitive to the parameters. However, variables such as hoop strain energy, are. Sensitivity to the initial state of compression is shown in Table 4. Sensitivity to the final state of compression is shown in Table 5.

TABLE 1. SAMPLE OUTPUT FROM THE COMPUTER PROGRAM: COMPACT

DIMENSIONS OF CYLINDRICAL SLEEVE SIGNITERNAL DIAMETER 3.2004 CM OUTSIDE DIAMETER 3.8100 CM EFFECTIVE HEIGHT 3.5560 CM	1.2600 1.5000	INCH	BODY		
PUNCH TRAVEL LIMIT 0.4157 CM		INCH			
PROPERTIES OF RDX , USED IN THE EXP	.osive fill:				
	0.0 MICRON				•
RADIAL GRID ELEMENT OF RDX	8.0 MICRON				
	06.0 SQ CM/ .820 G/CC	G			
	300 CAL/G/	ነበፍር የ			
	.00000E=04		S/DEG K		
	.28205E-03				
	.80000E+10				
	2.20000E-01				
	.07143E+10				
	. 985				
THERMAL PROPERTIES OF GRENADE MATER	'AIQ.				•
					•
	3.00100E-02				
	1.78917E+00		EO K		
	3.01485E-01 5.95200E-04				
	1.28882E-03				
	62000E-02				
	7.87000E+00				
	.25000E-01		EG K		
	1.65000E-01				
THERMAL DIFFUSIVITY OF STEEL +	1.67726E-01	SQ CM/S	}		
			 		_
INITIAL CONDITIONS AT START OF FINA		TION:			
	CC PSI	35.195	MPA		
HEIGHT OF HE COLUMN 2.58468 C		320 132	PIFA		
	esi	203.21	MPA		
HOOP STRAIN IN BODY +9.82452E-					
HOOP STRAIN ENERGY 1.1795 J		0.2819	CAL		
	OULE	0.1166			
CUM WORK BY PUNCH 1.6675 J	OULE	0.3985	CAL		
SIMULATED CONDITIONS DURING HE CONS	OLIDATION				
PUNCH HE HE MAX HOO		PUNCH	THERM	TEMP	CYCLE
DISPL DENS PRESS STRESS STR		WORK	WORK	RISE	TIME
(CM) (G/CC) (KPSI) (KPSI) (MU	U) (CAL)	(CAL)	(CAL)	(D K)	(MS)
0.0010 1.4406 5.115 29.53 98	1.5 0.283	0.467	0.067	0.007	2.75
	_				

TABLE 2. SENSITIVITY OF SIMULATION OUTPUT TO BODY WALL THICKNESS

Output	Body Wa	ll Thickness	(inch)
Variable	0.110	0.115	0.120
Loading Time (ms)	266	268	269
Punch Travel (mm)	3.976	4.031	4.087
Punch Work (cal)	57.5	57.3	57.1
Max Hoop Stress (kpsi)	159.1	151.8	144.9
Hoop Strain Energy (cal)	7.765	7.724	6.815

TABLE 3. SENSITIVITY OF SIMULATION OUTPUT TO EFFECTIVE BODY LENGTH

Outmat	Effective Leng	gth of Body (inch)
Output Variable	1.0	1.4
Loading Time (ms)	268	268
Punch Travel (mm)	4.031	4.031
Punch Work (cal)	57.3	57.3
Max Hoop Stress (kpsi)	151.8	151.8
Hoop Strain Energy (cal)	5.196	7.274

TABLE 4. SENSITIVITY OF SIMULATION OUTPUT TO INITIAL SYSTEM STATE Internal Diameter (cm) 3.226 Effective Length (cm) 3.556

Out	Initial Punch Pressure/HE Density			
Output Variable	4.75/1.42	(kpsi)/(g/cc)	5.10/1.44	
Loading Time (ms)	281		268	
Punch Travel (mm)	4.392		4.031	
Punch Work (cal)	59.6		57.3	
Max Hoop Stress (kpsi)	151.8		151.8	
Hoop Strain Energy (cal)	7.27		7.27	
Remain Heat Energy (cal)	49.5		47.2	

TABLE 5. SENSITIVITY OF SIMULATION OUTPUT TO PEAK PUNCH PRESSURE

Out sut	Peak Punch I	D		
Output Variable	25 (kpsi)	20 (kpsi)	Percent Decrease	
Max HE Density (g/cc)	1.697	1.678	1.1 %	
Final HE Density (g/cc)	1.676	1.661		
Max Hoop Stress (kpsi)	144.8	115.8	20.0 % 36.0 %	
Hoop Strain Energy (cal)	6.80	4.35		

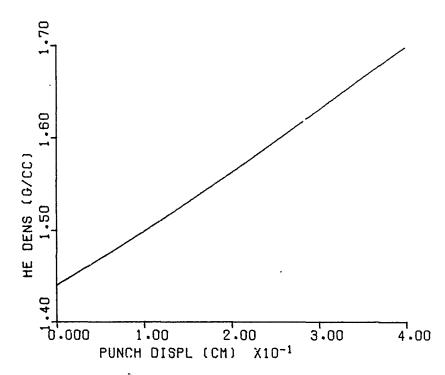


Figure 1. Avg Density of HE as a Function of Punch Displacement

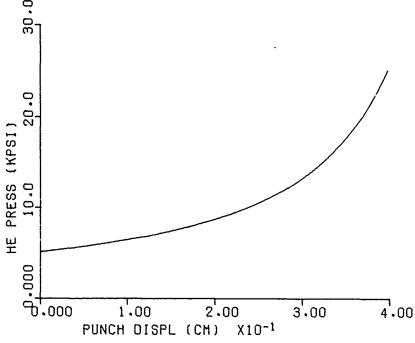


Figure 2. Max Pressure on the HE Versus Punch Displacement

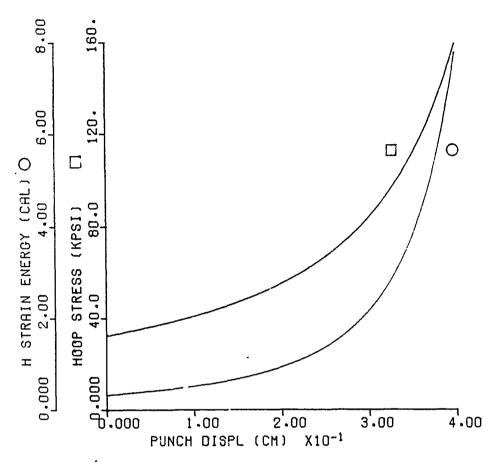


Figure 3. Hoop Strain Energy and Max Hoop Stress in the Grenade Body Versus Punch Displacement

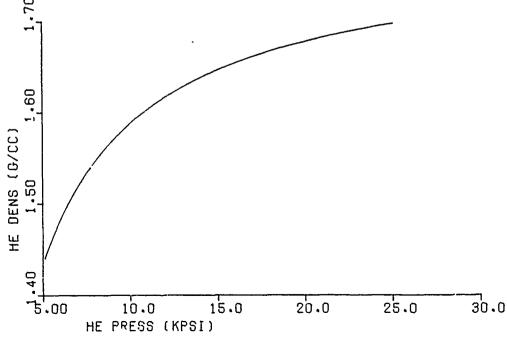


Figure 4. Avg Density of the HE Versus Max HE Pressure

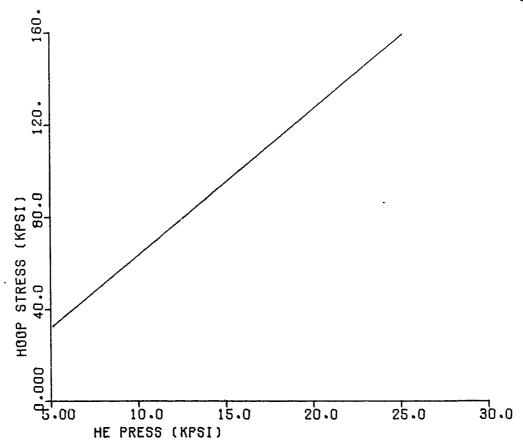


Figure 5. Max Hoop Stress in Grenade Body Versus Max HE Pressure

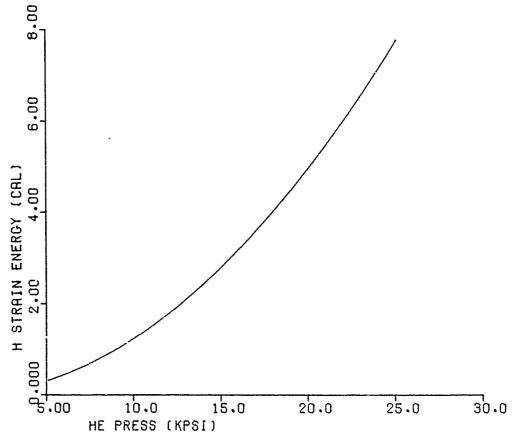


Figure 6. Hoop Strain Energy in the Grenade Body Versus Max HE Pressure

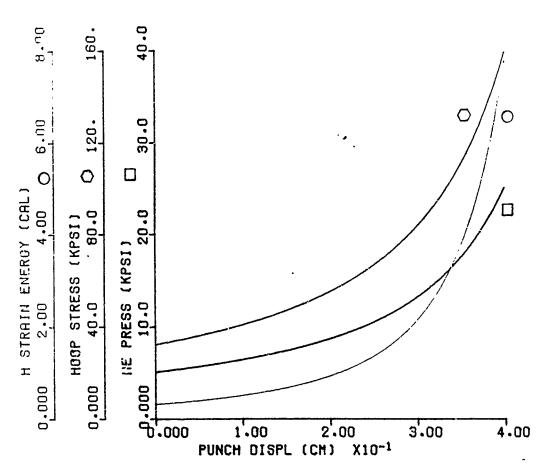


Figure 7. Comparison of Several Variables as Functions of Punch Displacement

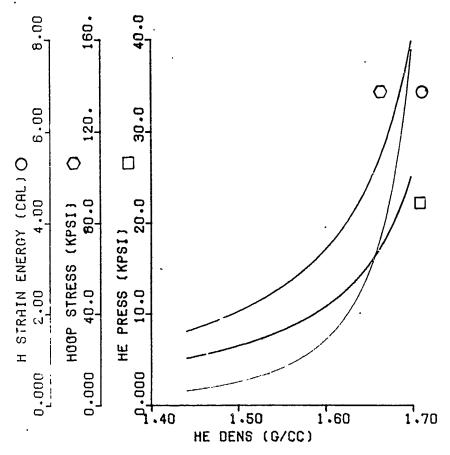


Figure 8. Comparison of Several Variables Versus Avg HE Density

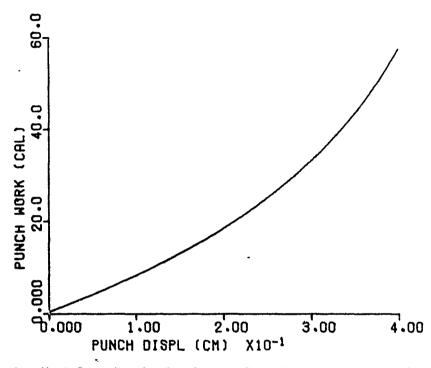


Figure 9. Work Done by the Punch as a Function of Punch Displacement

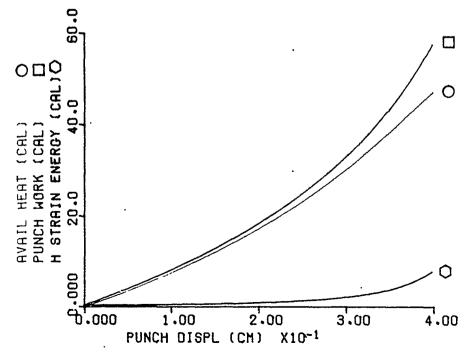


Figure 10. Comparison of Energy Components as Functions of Punch Displacement

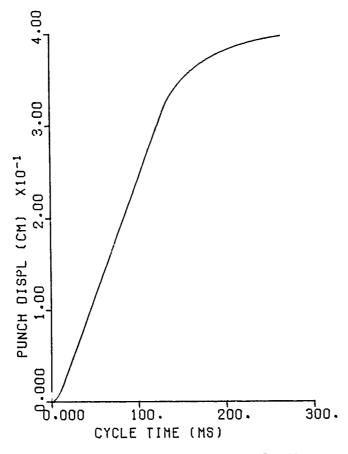


Figure 11. Punch Displacement Versus Cycle Time

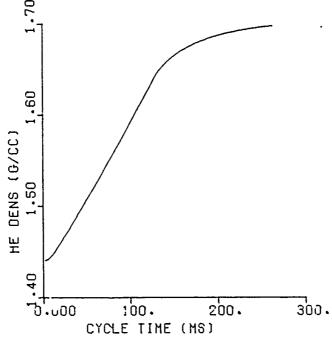


Figure 12. Avg Density of the HE Versus Cycle Time

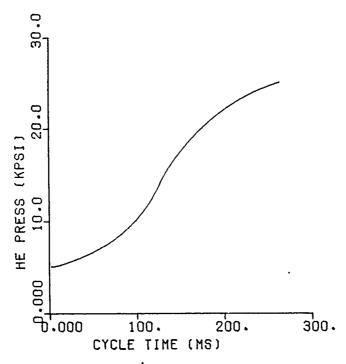


Figure 13. Max HE Pressure Versus Cycle Time

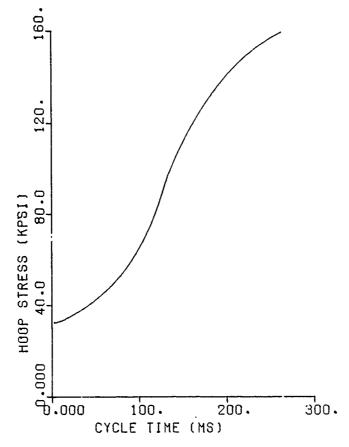


Figure 14. Max Hoop Stress Versus Cycle Time

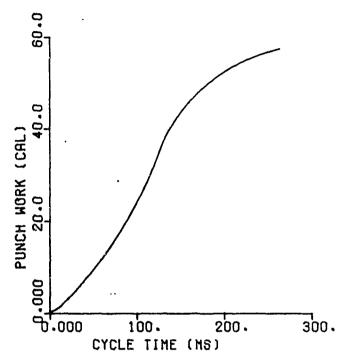


Figure 15. Punch Work Versus Cycle Time

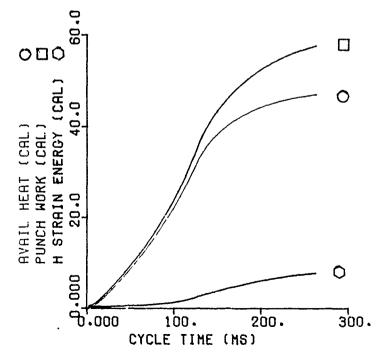


Figure 16. Comparison of Energy Components as Functions of Cycle Time

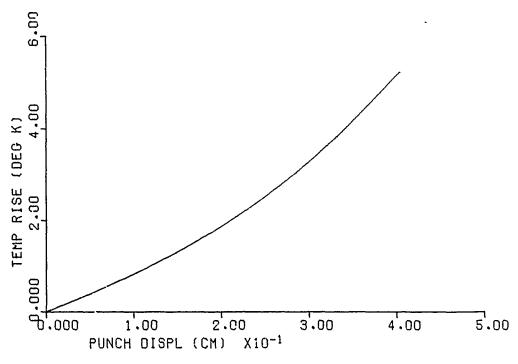


Figure 17. Avg HE Temperature Rise as a Function of Punch Displacement

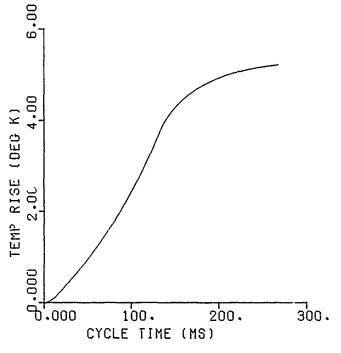


Figure 18. Avg HE Temperature Rise as a Function of Cycle Time

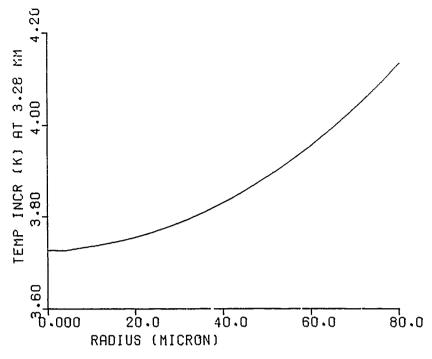


Figure 19, Radial Profile of Temperature in an RDX Particle (at 3.28mm Punch Travel)

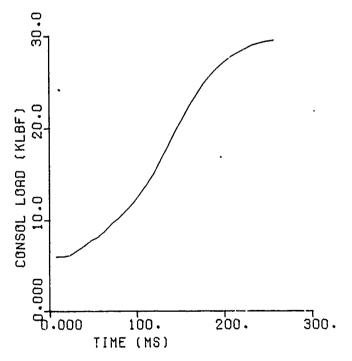


Figure 20. Experimental Consolidation Load Versus Cycle Time

10. Discussion of Results

For the most part the graphical results do not require comment. In some cases, however, my observations may help to clarify an issue or to make an Interpretation. The first ten figures show important variables as functions of punch displacement (or travel), whereas the last ten are largely functions of cycle time. By "cycle time" is meant time, measured from the start of the impression phase of the final consolidation cycle. Another semantic point: Maximum pressure on the HE fill is identical to the punch pressure. These terms are used interchangeably. In Figure 1 the average density of the HE is shown plotted as a function or punch displacement. Altho this relationship appears linear, it is not quite. Departure from linearity is due to progressive hoop strain in the body wall as the punch advances. Figure 2 shows the punch pressure as a function of punch travel. It is clear that the most rapid increase in pressure happens during the final half of the forward displacement. This functional form indicates that the HE behaves compressionally as a very nonlinear spring. Hoop strain energy and maximum hoop stress in the grenade body are shown as functions of punch displacement in Figure 3. It is noted that strain energy is a quadratic function of stress. Also observe that hoop stress builds rapidly near the end of forward travel. of the punch. For these reasons the hoop strain energy in the body is strongly dependent upon punch travel during the last quarter of the forward stroke. Therefore, one expects that brittle fracture would most probably occur in the last quarter of compressional punch travel. Only in this last quarter does the body strain energy exceed about two calories. Note that the max strain energy is nearly 8 calories. Figure 4 is a cross-plot of avg HE density versus punch pressure. It is apparent that at about 15 kpsi pressure, the rate of increase of HE density diminishes rapidly. Max body hoop stress is shown as a function of punch pressure in Figure 5. Maximum hoop stress occurs at the inside surface of the body wall. As predicted by this model, max hoop stress is a linear function of punch pressure. Figure 6 shows the hoop strain energy in the grenade body as a function of punch pressure. When plotted in this manner, hoop strain energy exhibits a more uniform rate of increase than when plotted versus punch displacement, as in Figure 3. In both Figures 7 and 8, hoop strain energy, max hoop stress, and max HE pressure are compared functionally. Whether the abscissa is punch travel (Fig 7) or avg HE density (Fig 8), it is seen that strain energy shows the greatest relative variation of these variables. Hoop stress and punch pressure exhibit more uniform growth. Figure 9 shows work done by the punch as a function of punch displacement. Because of the greater stiffness of the HE fill near the end of forward travel of the punch, the rate of work done by the punch increases progressively. This work is converted into both elastic and inelastic energy components. The elastic components (a) hoop strain energy in the grenade body, and (b) bulk-compression strain energy of RDX crystals in the HE. Punch work, available heat, and hoop strain energy are shown as functions of punch travel in Figure 10. The work performed by the punch is the total available energy. This is divided among the components: (a) body hoop surain energy, (b) RDX strain energy, (c) thermal loss, and (d) available thermal energy. Of these components, only the largest two are shown here.

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predicted rise in HE temperature, when loading to 25.4 kpsi, is 5.2 deg K. This compares with a measured value of 3.9 deg K, when loading to the same point (max load of 33 klbf) *. A max external hoop stress is reported (Ref 2b) as 105.5 kpsi for a peak punch load of 26.5 klbf. Simulated external stresses were obtained for the me max compression (21.2 kpsi) for two values of body ** wall thickness. For a 0.12-inch wall, the predicted peak, external hoop stress is about 102 kpsi; whereas, for a 0.11-inch wall, this value is about 107 kpsi. Agreement between the simulated and measured values of these variables is surprisingly close. With respect to the third variable used for comparison, one should note that functions, not scalars, are being compared. A single, arbitrary parameter in the simulation--max punch speed--is selected to obtain a good match of theory and experiment. This parameter adjusts the kinematics of the simulation to that of the experiment; however, this parameter does not affect the shape or the amplitude of the load function. Therefore, good agreement in the simulated and measured load functions constitutes a challenging test of the validity of the model. The experimental load-time curve, shown in Figure 1 in Ref 2i, is matched by simulated results well within experimental error (+ or - 1.6 klbf) over the loading interval of about 260 ms from start of cycle. The experimental load function is shown for comparison in Figure 20, for just the compression phase of the load cycle. No systematic difference is apparent between model and measured functions.

13. Methodology

Before discussing math models for various phenomena, I will outline the procedure followed in simulating the loading process. The independent variable in the simulation is the displacement (or travel) of the punch. The kinematics of the process are obtained thru the functional relationship between punch displacement, xp, and punch speed, sp. Only the compression-, or loading, -phase of the consolidation

Temperature in this environment is very difficult to measure. Ref 2b states that two types of transducers were used -- theromcouple and nickel-foil gage. The nickel-foil gages were too fragile: three of four were damaged. Two measurements, at different locations within the HE, were made with thermocouples. One measurement was made just above the tip of the punch, with the HE consolidated at at max load of 43 klbf. The reported temperature r_se in this case is 18.9 deg K. This value seems much too high; and, in any case, does not represent a volume-average temp rise. The other reported measurement was made between the axis and wall near the top or cap end of the cavity. This is the value (3.9 deg K) cited here, which was obtained at the lowest reported peak load (33 klbf). Parenthetically, one observes that if the temperature variation within the HE is a great as measured, this fact would support the assertion that considerable differential flow and, possibly, grain fracture occurs within the HE.

^{**} Recall that the grenade body is represented by a uniform cylindrical sleeve. Wall thickness of the sleeve is the parameter in this instance. The parametric variation studied (0.010 incn) is about three times the body-to-body standard deviation for a particular body producer.

cycle is simulated. After initialization of the grenade to a given HE density and corresponding punch pressure, the following steps are followed until the desired terminal state is achieved:

- (a) The punch is advanced thru a small increment. This increases the density of the HE. (One micrometer is the increment used here.)
- (b) The max pressure on the punch and the pressure gradient in the HE are calculated at the current average HE density. The incremental work performed by the punch, to advance incrementally at this pressure, is added to cumulative punch work.
- (c) The (one-dimensional) hoop strain in the cylindrical sleeve, which represents the grenade body, is calculated. Max strain is reached at the inside radius. Elastic behavior of the metal is assumed.
- (d) Since the hoop strain increases the HE cross-section, calculate the additional displacement of the punch (at this pressure) required to preserve HE density at the value calculated previously with no incremental body strain.
- (e) At each movement of the punch, by delxp, add the punch work to the cumulative punch work. The isobaric, incremental work is just the total force times delxp, where the total force is the product of the current values of HE cross-section and punch pressure. Updates of punch work are required at steps (b) and (d).
- (f) Since the RDX particles within the HE experience elastic volumetric compression, calculate the strain and the strain energy in the RDX mass.
- (g) To obtain the total available heat energy, subtract the sum of hoop-strain energy plus RDX strain energy plus thermal loss to body from the punch work. (All cumulative energy terms are calculated in joules and reported in units of J and calories.)
- (h) The total incremental punch displacement (since prior loop pass) has consumed a time increment which is the ratio of the displacement increment to the punch speed. This speed is obtained as a function of current value of xp. Time is updated by addition of the time increment.
- (i) The time increment is also used to calculate the thermal flux to the RDX particles. Flux is the incremental heat/total particle surface/time increment. At the program user's option, the temperature distribution as a function of RDX particle radius is obtained by numerically solving the diffusion equation, using the average thermal flux as the outside boundary condition.
- (j) The heat generated incrementally during this loop pass is used to calculate the volume-average temperature rise of the HE, by dividing by the heat capacity of the HE. HE temperature is updated by this increment.

- (k) The incremental heat loss from the HE to steel body is calculated using current average values of HE temperature and steel temperature. Continuity of temperature and thermal flux at the HE-wall boundary are assumed. The incremental heat loss is used to update the total heat loss and the average steel temperature.
- (1) Current values of the state variables are printed optionally at multiples of the number of loops. Regardless of option, saved values—such as punch displacement and available heat—are stored.
- (m) If the required terminal state has not been reached, loop back to (a) for another pass. Otherwise, step and print final results. Among the final results are: volumetric increase in the HE due to elastic recovery and the associated decrease in HE density. (If a portion of the mass of RDX experienced grain fracture, some of the RDX strain energy would be thermalized, and the volumetric recovery would be smaller than calculated.)

14. Model Equations

The average density of the HE is, by definition, the ratio of HE mass, M, to volume, V. But, the volume of HE depends upon the cross section and length (or height) of the HE column. Denoting cross sectional area by A, initial column length by xo, and punch displacement by xp, the volume of the HE is given by

$$V = A(xo - xp), 0 le xp le xpmax. (1)$$

Denoting the average (bulk HE) density by E(rho),

$$E(rho) = M/(A(xo - xp)).$$
 (2)

Altho not explicitly written as a function of displacement, the area A should be considered a function of xp, since the max hoop strain, eps, is a function of xp, and since

$$A = Ao(1 + eps), \qquad (3)$$

where Ao is the unstrained, internal cross section of the body. A formula for E(rho) is derived in Ref 2d in terms of the theoretical maximum density of the HE, TMD, and of the punch pressure, p. This formula is repeated here:

$$E(rho)/TMD = 1 + (0.8/c)ln((p+0.8-c)/(p+0.8)),$$
 (4)

where c is the difference between maximum and minimum pressure within the volume of the HE. This result assumes a constant pressure gradient within the volume. The basis of the derivation is a constituative relationship between local HE density, rho, and local pressure, p:

$$rho/TMD = p/(p + 0.8),$$
 (5)

where the constant 0.8 is given in kpsi.

Based on several experimental studies, cited in Ref 2d, the parameter c, which is a measure of the pressure gradient, is found to be directly proportional to the maximum pressure on the HE. Thus, parameter

$$\mathbf{c} = \mathbf{g} \, \mathbf{p} \,, \tag{6}$$

where g is a constant. From Ref 2e, the value of g which best fits data on average bulk density is 0.6. This value is used here. (In earlier experiments at KSAAP and MAAAP, a value of 0.8 for g was indicated.)

15. Equation (4) for average density is not readily inverted in order to give punch pressure as a function of E(rho). This fact poses no problem numerically, however, since the relationship of these variables is monotonic. In a subroutine of COMPACT, punch pressure is calculated as a function of E(rho) by an iterative Newton-Raphson method. Given the punch pressure from (4), one can calculate the maximum hoop (i.e., tangential) stress within the wall of the cylindrical sleeve, representing the steel body, by Lame's formula (Ref 2h). The hoop stress at radius r is s(r):

where p is the pressure on the inside wall and po is the pressure on the outside of the wall. The internal radius of the sleeve is a and the outside radius b. The radius to an arbitrary point is r. In the present application, the maximum internal pressure is identified with the punch pressure, and the outside pressure is taken as atmospheric pressure, i.e., the outside of the wall is assumed unsupported. The maximum stress occurs at the inner wall, where r = a. The max body stress is, then,

If the body remains elastic thruout, the maximum strain is simply

$$eps = s(a)/Ey , (9)$$

where Ey is Young's modulus for steel. As indicated above, the cross section of the HE is related to xp functionally thru equations (2), (3), (4), (8), and (9).

16. The total hoop strain energy in the body is obtained by integrating the strain energy per unit volume,

over the volume of the cylinder wall subject to strain. Let the effective cylinder height experiencing this hoop strain be denoted by h. Then, using equation (7), the hoop strain energy is given by:

HSE =
$$(pi h/Ey)(T + T + T)$$
, (10)

where the indexed terms are functions of the dimensions a and b and of the punch pressure p.

where the auxiliary factors A and B are given by

$$2 2 2 2 2$$

$$A = (a b (p - po))/(b - a)$$

$$2 2 2 2 2$$

$$B = (p a - po b)/(b - a).$$
(11b)

Energy is also stored in the RDX crystals as elastic strain. If a local hydrostatic condition is assumed for each RDX particle, an expression for the total RDX strain energy can be derived by integrating stain energy per unit volume over the total volume. Let the bulk modulus of RDX be denoted by Bm. Also, let the pressure at a normalized axial coordinate x be denoted by p(x). The assumption of a constant pressure gradient in the HE means that

$$p(x) = p(1 - gx),$$
 0 le x le 1, (12)

where, as above, p is the punch pressure and g is a constant. The RD λ strain energy per unit volume, at x, is

By integrating this expression, with p(x) given in (12), over the total volume of RDX, Vr, one obtains the RDX strain energy, RSE:

$$2 2 RSE = (Vr p / (2 Bm))(1 - g + g / 3) . (13)$$

Work performed by the punch is, of course, the source of both elastic and inelastic energy components. The punch work, W, is the integral over punch travel of

A p
$$delta(xp)$$
, (14)

where delta(xp) is the incremental displacement of the punch, given the constant punch pressure p and HE cross section A. For 3-digit accuracy the value of delta(xp) must be quite small--about 5 microns or less. At any instant, the total heat residing in the HE is the work performed by the punch minus (elastic) potential energy, of body and RDX, and minus heat energy lost to the body. The heat loss is relatively small, but is included to be complete. Denoting the available thermal energy in the HE by H and the heat lost by L, an energy balance requires that

$$H = W - HSE - RSE - L . (15)$$

The average temperature rise, avgUh, of the HE, with heat capacity Ch is, then,

$$avgUh = H/Ch. (16)$$

Similarly, the average temperature rise of the steel body, avgUs, is obtained from L and the heat capacity of the body, Cs:

$$avgUs = L/Cs . (17)$$

17. A mathematically exact calculation of the heat loss from HE to body wall requires the solution of the diffusion equations in both HE and steel wall. The diffusion equation in the HE involves a heat source term, representing the rate of energy production per unit mass. To describe this source term requires an assumption concerning HE flow within the cavity, since viscous effects produce the heat. But, details of the HE flow are not modeled here. To escape this dilemma, I have simply assumed that the source term is independent of position (but, not of time). (Actually, this isotropic assumption implies something concerning HE flow, but is not pursued.) Denoting the rate of heat released per unit mass per unit time as Q, one can write

$$Q = delta(H)/delta(t)/M, (18)$$

where delta(H) is the increment of heat added during the time increment delta(t). Notationally, let U(r) be the temperature at radial position r. When r is less than or equal to the inner radius, a, the temperature pertains to the HE. For r between a and outside radius, b, the temperature describes the steel wall. Further, let a terminal "1" indicate a property of the HE, and a terminal "2" indicate a property of the steel. Thus, alphal denotes the thermal diffusivity of HE, Cl denotes the specific heat of HE, and Kl denotes the thermal conductivity of HE. (The particular HE is Comp A-5.) Partial differentiation by r and by t are denoted, respectively, by subscripts r and t. With these conventions, the temperature in the HE is given by

$$U = (alpha1/r)(r U) + Q/C1, 0 < r < a.$$
 (19)

The diffusion equation for the cylindrical wall is:

$$U = (alpha2/r)(rU), \qquad a < r < b. \qquad (20)$$

Boundary conditions are needed at radial positions: 0, a, and b. At r=0, cylindrical symmetry requires that no heat is transported across the center. Thus,

$$(U(0)) = 0. (21a)$$

At r = a, temperature is the same on each side of the boundry; and, the thermal flux out of the HE must equal flux into the steel wall.

$$U(a-) = U(a+)$$

and

$$K1 (U(a-)) = K2 (U(a+)),$$
 (21b)

where a- and a+ means r approaching a from below and from above, resp. Finally, the boundary at r = b can be treated as thermally insulated, since still air outside the cylinder is a good insulator.

$$(U(b)) = 0. (21c)$$

The rate of heat loss from HE is obtained from the following equation:

$$delta(L)/delta(t) = -K1 Awall (U(a-)),$$
(22)

where Awall is the area of HE contacting the wall. The cumulative heat loss, L, is obtained by numerical integration of the time derivative in (22). A numerical solution of (19) thru (22) is obtained by a integrating a set of total differential equations for U(i), defined on a radial grid r(i), i=1,2,.... These equations are derived from (19,20,21) by a conventional forward-difference scheme. The implementing computer code is located on page A-7 of Annex A.

18. An Approximation for Heat Loss

Because the importance to this simulation of the heat loss from the HE is relatively small, the approach taken in paragraph 17 may entail too much computation. An approximation is given here which, tho not as accurate, may suffice. The approximate rate of loss makes use of the great difference in the thermal diffusivities of HE and steel. Because RDX has a very low diffusivity, heat loss is not felt very far into the HE from the cylinder wall. Therefore, a very steep temperature gradient exists in the HE near the wall. However, since heat diffuses rapidly in steel, a rather shallow radial temperature gradient gradient exists in the wall. Thus as a first approximation, the gradient in the steel is taken as zero, which implies that the temperature at the wall is nearly avgus. Additionally, assume that the temperature in the HE drops linearly from avguh to the value at the wall over a radial interval DELR. The negative gradient in the HE near the wall is approximated by

The approximate rate of heat loss from the HE is, from (22), the product of the factors Awall, K1, and the negative gradient in (23). The increment DELR is effectively the distance traveled internally from the wall at which the temperature rises linearly to avgUh. A natural length for measuring temperature change is the diffusion length, the distance a temperature spike will propagate by diffusion in time t. The diffusion length in the HE is

Empirically, a good approximation for the temperature gradient at the wall, during the compression phase of loading, is obtained when DELR is 4 times the diffusion length.

19. Temperature in an RDX Particle

Among the physical processes occurring during loading, thermal diffusion in RDX particles appears to be of minor importance. If the heating rate of the surface of a typical RDX particle were sufficiently great, thermally (as well as mechanically) induced stresses might play a role in grain fracture. Actually, now this does not appear to be the case; but, this idea led to a study of the temperature profile with respect to radius in an RDX particle. The mean RDX particle diameter is about 160 microns for Class 1 RDX, which is presently being used in Comp A-5. Therefore, a spherical particle of this diameter was selected for the temperature study. The RDX specific surface—area per unit mass—was calculated for a uniform granulation of this diameter. Denoting the specific surface of the RDX by spS, the heat flux directed at a particle is given by

$$q = spS Q (Mass RDX/Mass HE)$$
, (25)

where the generation rate per unit mass of HE, Q, is given by (18). Notationally, let the RDX particle diameter be d, and let a terminal "3" denote a material property of RDX. Thus, K3 denotes the thermal conductivity of RDX. Additionally, denote the RDX temperature rise at particle radius r by U(r). With this notation and subscript conventions, one can write the boundary condition at the surface of the typical particle as

$$q = -K3 (U(d/2))$$
, (26)

assuming that all the flux is absorbed. By spherical symmetry, the boundary condition at particle center is:

These boundary conditions are applied to the spherically symmetric form of the diffusion equation in a particle. Functional dependence of U upon r is suppressed here:

$$U = \text{alpha3} ((2/r)U + U) , 0 < r < d/2. (28)$$

Solution of the RDX diffusion equations is an option in the program COMPACT. The numerical method uses a radial grid of points at which the temperature is evaluated. The differential equations on this grid are developed from (28) using a central-difference approximation to spatial derivatives. An update is performed using two half-steps at each step in time. Equations are found on page A-8 in Annex A. Quite small time steps, delta(t), are required for numerical stability of this method. To insure that stability is met, the incremental punch travel, delxp, is kept small (1 micron), since this and punch speed, implicitly determine delta(t). Additionally, when calculating RDX particle temperature, the initial punch speed is taken to be its maximum value, (with a slight loss of fidelity to kinematics).

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ANNEX A

SOURCE PROGRAM FOR SIMULATING THE STATE OF AN M42/M46 GRENADE DURING PRESS LCADING

PREAMBLE ''COMPACT NORMALLY MODE IS REAL DEFINE PRESS AS A REAL FUNCTION GIVEN 2 ARGUMENTS DEFINE RHOTRANS AS A REAL FUNCTION GIVEN 2 ARGUMENTS DEFINE STRN. ENER AS A REAL FUNCTION GIVEN 2 ARGUMENTS DEFINE ESTRN.RDX AS A REAL FUNCTION GIVEN 1 ARGUMENT 6 DEFINE RPSPEED AS A REAL FUNCTION GIVEN 1 ARGUMENT DEFINE FLAGU AS AN INTEGER VARIABLE DEFINE EYOUNG, LBODY, RINT, ROUT, GRAD. PRS, XPMAX AS REAL VARIABLES DEFINE BM.RDX, EY.RDX, PR.RDX, VOL.RDX AS REAL VARIABLES 10 11 END ''PREAMBLE MAIN ''COMPACT 1 2 ''PROGRAM SIMULATES THE COMPACTION OF THE HE AND THE STRAINING OF THE ''GRENADE BODY DURING PRESS LOADING OF AN M42 GRENADE. A BODY IS REP-5 "RESENTED AS A CYLINDRICAL SLEEVE OF CONSTANT HEIGHT. ''PARAMETERS: MASS OF THE HE (G). 8 ''MASSHE CROSS-SECTIONAL AREA (SQ CM) OF THE HE DURING PRESSING. 9 ''AXHE 10 ''AXHEO INITIAL CROSS-SECTION OF THE HE (SQ CM) ''RINT INITIAL INTERNAL RADIUS OF THE GRENADE (CM). 11 INITIAL OUTSIDE PADIUS OF THE GRENADE (CM). ''ROUT 12 EFFECTIVE LENGTH OF THE BODY SUBJECT TO HOOP STRAIN (CM). 13 ''LBODY 14 ''X0 INITIAL EFFECTIVE HEIGHT OF THE HE CYLINDER (CM). 15 ''XP DISPLACEMENT OF THE PUNCH (CM). ''XPMAX MAX DISPLACEMENT OF THE PUNCH (CM). 16 ''DELXP INCREMENT IN DISPLACEMENT OF THE PUNCH (CM). 17 18 ''RHOHEO DENSITY OF THE HE PRIOR TO FINAL CONSOLIDATION (G/CC). AVG DENSITY OF THE HE DURING CONSULIDATION (G/CC). 19 ''RHOHE 20 ''TMDHE THEORETICAL MAX DENSITY OF THE HE (G/CC). PRESSURE ON THE HE (KPSI) DURING CONSOLIDATION. 21 ''PHE INTERNAL PRESSURE ON THE HE BODY (PASCAL). 22 ''PINT OUTSIDE PRESSURE ON THE BODY (PA). 23 ''POUT 24 ''MAXKPSI MAX PRESSURE (KPSI) TO STOP THE SIMULATION. 25 ''EYOUNG YOUNG'S MODULUS OF THE BODY STEEL (PA). ''DENS.STEEL DENSITY OF STEEL IN BODY (G/CC). 26 ''SPHT.STEEL SPECIFIC HEAT OF STEEL (CAL/G/DEG K). 27 28 ''TCOND.STEEL THERMAL CONDUCTIVITY OF STEEL (CAL/CM/S/DEG K). 29 ''DIFUZ.STEEL THERMAL DIFFUSIVITY OF STEEL (SQ CM/S). 30 'STGPA BODY HOOP STRESS (PA). ''SIGKPSI BODY HOOP STRESS (KPSI). 31 ''HSTRAIN BODY HOOP STRAIN. 32 33 ''GRAD.PRS COEFFICIENT FOR THE PRESSURE GRADIENT WITHIN THE HE. 34 ''MASS.RDX MASS OF RDX IN HE (G). 35 ''FRACT.RDX MASS FRACTION OF RDX IN HE. 36 ''VOL.RDX VOLUME OF RDX CRYSTALLS (CC). 37 ''BM.RDX BULK MODULUS OF RDX (PA). 38 ''EY.RDX YOUNG'S MODULUS FOR RDX (PA). 39 ''PR. RDX POISSON'S RATIO FOR RDX. 4C ''SPS.RDX SPECIFIC SURFACE FOR RDX (SQ CM/G).

```
THEOPETICAL MAX DENSITY OF RDX (G/CC).
  41
       ''TMD.RDX
  42
       ''SPHT.RDX
                     SPECIFIC HEAT OF RDX (CAL/G/DEG K).
  43
       ''TCOND. RDX
                     THERMAL CONDUCTIVITY OF RDX (CAL/CM/S/DEG K).
       ''DIFUZ.RDX -
  44
                     THERMAL DIFFUSIVITY OF RDX (SQ CM/S)
       "MASS.STER _
  45
                     MASS STERIC ACID (G).
  46
       ''DENS.STER DENSITY OF STERIC ACID (G).
  47
       ''SPHT.STER
                     SPECIFIC HEAT OF STERIC ACID (CAL/G/DEG K).
  48
       ''HSE
                     BODY (ELASTIC) HOOP-STRAIN ENERGY (JOULE).
       ''EE. RDX
  49
                     ELASTIC ENERGY IN THE RDX CRYSTALS (JOULE)
  50
       ''HENERGY
                     (THERMAL) ENERGY AVAILABLE FOR HEATING THE RDX.
  51
                     SPACE INTERVAL IN RDX-PARTICLE RADIAL GRID (CM).
  52
       ''DELTAT
                     TIME STEP IN INTEGRATING THE "EAT THANSFER DIFF EQNS.
  53
      DEFINE FLAGUHE, I, J, K, KPRINT, KUPRINT, L, M, N, NGRID, NHE, NST AS INTEGER
       VARIABLES
      DEFINE ANSWER AS A TEXT VARIABLE
  56
      DEFINE UV, UOV, U1V AS REAL, 1-DIMENSIONAL ARRAYS ''RDX PARTICLE TEMP
      DEFINE JHEV, UHEOV, USTV, USTOV AS REAL, 1-DIMENSIONAL ARRAYS ''HE TEMP
  57
  58
      ''ASSIGN CONSTANTS.
  59
   60
  61
           LET FLAGUHE=0 ''IS 1 TO FLAG CALC OF RADIAL DIST OF UHE
           LET NGRID=11
  62
  63
           LET NHE=21
  64
           LET NST=5
   65
           LET DELTAR=0.0008 ''CM STEP IN RDX XTAL (8 MICRON)
  66
           RESERVE UV(*).UOV(*).U1V(*) AS NGRID
  67
           RESERVE UHEV(*), UHEOV(*) AS NHE
  68
           RESERVE USTV(#), USTOV(#) AS NST
  69
           FOR I=1 TO NHE, LET UHEOV(I)=0.0
   70
           FOR I=1 TO NST, LET USTOV(I)=0.0
           FOR I=1 TO NGRID DO
   71
               LET UV(I)=0.0
   72
               LET UOV(I)=0.0
   73
   74
               LET U1Y(I)=0.0
   75
           LOOP ''TO INITIALIZE THE RDX TEMPERATURE PROFILE
           LET HTLOSS=0.0
   76
   77
           LET KPRINT=10
           LET KUPRINT=100
   78
   79
           LET DELXP=0.0001
   80
           LET GRAD. PRS=0.6
   81
           LET PA.KPSI=6.89474*10**6 ''PASCALS PER KPSI CONVERSION
   82
           LET CALPJ=0.239 ''CALORIES PER JOULE CONVERSION
  83
           LET EYOUNG =2.06843*10**11 ''PA. YOUNG'S MODULUS FOR STEEL
  84
           LET DENS, STEEL=7.87 ''G/CC
   85
           LET SPHT.STEEL=0.125 ''CAL/G/DEG K
   86
           LET TCOND.STEEL=0.165 ''CAL/CM/S/DEG K
   87
           LET DIFUZ.STEEL=TCOND.STEEL/DENS.STEEL/SPHT.STEEL
   88
           LET LBODY=3.556
           LET RINT=1.6256 ''FOR 110 MIL WALL THICKNESS
   89
           LET RINT=1.6002 FOR 120 MIL WALL THICKNESS
   90
   91
           LET RINT=1.6130
   92
           PRINT 2 LINES WITH RINT
           THUS
   93
CURRENT VALUE OF INTERNAL RADIUS IS #. **** CM, FOR A 110 MIL WALL THICKNESS.
INPUT THE VALUE OF RINT WANTED. NOTE: 1.6002 (120 MIL WALL THICKNESS).
  96
           READ RINT
   97
           LET ROUT=1.9050 ''CM
```

```
98
         LET DRHE=RINT/(NHE-1)
 99
         LET DRST=(ROUT-RINT)/(NST-1)
100
         LET POUT=1.01353*10**5 ''PA (14.7 PSI)
101
         PRINT 1 LINE THUS
  INPUT THE MAX PUNCH PRESSURE. IN KPSI. TO STOP SIMULATION.
103
         READ MAXKPSI
104
         LET AXHEO=PI.C*RINT**2
105
         LET AXHE=AXHEO ''INITIALLY
106
         LET MASSHE=30.0 ''G
107
         LET FRACT.RD := 0.985 ''FRACTION OF RDX IN HE
         LET MASS.RDX=MASSHE*FRACT.RDX
108
109
         LET MASS.STER=MASSHE-MASS.RDX
110
         LET SPHT.STER=0.399
111
         LET DENS.STER=0.847
         LET TCOND.STER=3.8/10##4
112
         LET HTCAP.STER=MASS.STER*SPHT.STER ''CAL/DEG K
113
         LET TMD.RDX=1.82 ''G/CC
114
         LET TMDHE=1.78 ''G/CC FOR COMP A-5
115
         LET VOL. RDX=MASS. RDX/TMD. RDX ''VOLUME (CC) OF RDX CRYSTALS
116
         LET RAD. RDX=(NGRID-1)*DELTAR
117
118
         LET NPARTICLES=VOL.RDX/PI.C/4.0*3.0/RAD.RDX**3
         LET SPS.RDX=NPARTICLES#4.0#PI.C#RAD.RDX##2/MASS.RDX
119
         LET SPHT.RDX=0.300 ''CAL/G/DEG K
120
         LET HTCAP. RDX=MASS. RDX#52HT. RDX
121
122
         LET TCOND.RDX=0.0007 ''CAL/CM/S/DEG K
123
         LET DIFUZ. RDX_TCOND. RDX/TMD. RDX/SPHT. RDX
124
         LET FRACT.STER=1.0-FRACT.RDX
125
         LET SPHT.HE=FRACT.RDX#SPHT.RDX+FRACT.STER#SPHT.STER
         LET TCOND. HE=FRACT. RDX*TCOND. RDX+FRACT. STER*TCOND. STER
126
127
         LET DENS.HE=1.0/(FhACT.RDX/TMD.RDX+FRACT.STER/DENS.STER)
128
         LET DIFUZ.HE=TCOND.HE/DENS.HE/SPHT.HE
         LET UCON. HE=DIFUZ. HE/DRHE/DRHE
129
130
         LET UCON.ST=DIFUZ.STEEL/DRST/DRST
         LET DTIME=0.27 ''(S) TIME FOR TEMP DIFFUSION
131
         LET DL.HE=SQRT.F(DIFUZ.HE*DTIME) ''(CM) DIFFUSION LENGTH IN HE
132
         LET DELR.HE=4.0*DL.HE ''(CM) FOR APPROX THERM GRAD IN HE AT WALL
133
134
         LET PTT = TCOND. HE#DRST/TCOND. STEEL/DRHE
135
         LET F
                  =1.0/(1.0+BETA)
         LET L
                 =DIFUZ.RDX/DELTAR**2
136
         LET UOCON=DELTAR/TCOND.RDX
137
138
         LET U1CON=DIFUZ.RDX/DELTAR
139
         LET EY.RDX=1.8*10**10 ''YOUNG'S CONSTANT FOR RDX (PA)
         LET PR.RDX=0.22 ''POISSON'S RATIO FOR RDX
140
141
         LET BM.RDX=EY.RDX/3.0/(1.0-2.0*PR.RDX) ''BULK MODULUS FOR RDX
142
143
     ''GET INITIAL HE DENSITY FROM THE TERMINAL.
144
145
         PRINT 1 LINE THUS
  INPUT THE INITIAL DENSITY OF THE HE.
                                         SUGGEST 1.440 G/CC.
147
         READ RHOHEO
148
         PRINT 1 LINE THUS
  DO YOU WANT THE TEMPERATURE PROFILE IN AN RDX PARTICLE? (Y OR N).
150
         READ ANSWER
151
         IF SUBSTR.F(ANSWER.1.1) = "Y"
152
             LET FLAGU=1
153
         OTHERWISE
154
             LET FLAGU=0
```

```
ALWAYS
 155
       , ,
 156
 157
       ''GET INITIAL HEIGHT OF HE CYLINDER.
 158
 159
           LET VOL. HEO = MASSHE/RHOHEO
 160
           LET XO=VOL. HEO/AXHEO
       1 1
 161
       ''CALCULATE INITIAL BODY STRAIN DUE TO PRECONSOLIDATION.
 162
 163
 164
           LET PHE=PRESS(TMDHE, RHOHEO)
 165
           LET PINT=PA.KPSI*PHE
 166
           LET PINTO=PINT
           LET EE.RDX=ESTRN.RDX(PINTO) ''RDX STRAIN ENERGY FOR PRECONSOL COND
 167
           CALL STRS.STRAIN (PINT, POUT) YIELDING SIGPA, HSTRAIN
 168
 169
           LET SIGKPSI=SIGPA/PA.KPSI
 170
           LET HSTRESS=SIGPA/10**6
 171
       ''GET CROSS-SECTION OF HE. ACCOUNTING FOR BODY STRAIN.
 172
  173
           LET AXHE=AXHEO#(1.0+ABS.F(HSTRAIN))##2
 174
  175
  176
       ''REEVALUATE HEIGHT OF HE COLUMN TO GIVE DESIRED DENSITY.
  177
           LET XO=MASSHE/RHOHEO/AXHE
  178
  179
       ''GET MAX HOOP STRAIN AND MAX DENSITY.
  180
  181
           CALL STRS.STRAIN (MAXKPSI*PA.KPSI, POUT) YIELDING MAXSIGPA, MAXSTRAIN
  182
           LET MAXRHO=RHOTRANS(TMDHE, MAXKPSI) + 0.0001
  183
  184
           LET XPMAX=X0-MASSHE/MAXRHO/AXHEO/(1.0+ABS.F(MAXSTRAIN)) **2 + 0.007
  135
           LET AWALL=2.0*PI.C*(RINT*XO+RINT**2) ''WALL AREA FOR HEAT LOSS
  18€
           LET HVOL.STEEL=AWALL*(ROUT-RINT) ''VOL OF STEEL HEATED BY HE LOSS
  187
           LET HTCAP.STEEL=HVOL.STEEL*DENS.STEEL*SPHT.STEEL
  188
           LET USTEEL=0.0 ''INITIALIZATION OF TEMPERATURE OF STEEL BODY
           IET HSE=STRN.ENER (PINT.POUT)
  189
  190
           LE1 PWORK=HSE+RE. RDX
           LET HENERCYO-O O ''INITIAL HEAT ENERGY DEFINED AS ZERO (REFERENCE)
  191
           LET TIME=0.0
  192
           LET XP=0.0
  193
           LET XPO=XP
  194
           LET HOLDXP=0.0
  195
           SKIP 2 LINES
  196
           PRINT 5 LINES WITH 2*RINT,2*RINT/2.54,2*ROUT,2*ROUT/2.54,LBODY,
  197
  198
           LBODY/2.54, XPMAX, XPMAX/2.54
           THUS
  199
DIMENSIONS OF CYLINDRICAL SLEEVE SIMULATING THE GRENADE BODY
INTERNAL DIAMETER
                                 CM
                                           * * * * *
                                                   INCH
OUTSIDE DIAMETER
                                 CM
                                                   INCH
                                 CM
EFFECTIVE HEIGHT
                                                   INCH
PUNCH TRAVEL LIMIT
                                                   INCH
                                 CM
  205
           SKIP 2 LINES
  206
           PRINT 15 LINES WITH 20000*RAD.RDX.10000*DELTAR.SPS.RDX.TMD.RDX.
           SPHT. RDX, TCOND. RDX, DIFUZ. RDX, EY. RDX, PR. RDX, BM. RDX, FRACT. RDX
  207
           THUS
  208
PROPERTIES OF RDX , USED IN THE EXPLOSIVE FILL:
```

```
DIAMETER OF RDX PARTICLE
                                           MICRON
RADIAL GRID ELEMENT OF RDX
                                           MICRON
SPECIFIC SURFACE OF RDX
                                           SQ CM/G
THEORETICAL MAX DENSITY OF RDX
                                           G/CC
SPECIFIC HEAT OF RDX
                                           CAL/G/DEG K
THERMAL CONDUCTIVITY OF RDX
                                                   CAL/CM/S/DEG K
THERMAL DIFFUSIVITY
                      OF RDX
                                                   SQ CM/S
YOUNG'S CONSTANT FOR RDX
                                                   PASCAL
POISSON'S RATIO FOR RDX
BULK MODULUS FOR RDX
                                                   PASCAL
MASS FRACTION OF RDX IN HE
           PRINT 15 LINES WITH DRHE, DENS. HE, SPHT. HE, TCOND. HE, DIFUZ. HE, DRST,
  224
  225
           DENS. STEEL, SPHT. STEEL, TCOND. STEEL, DIFUZ. STEEL
  226
THERMAL PROPERTIES OF GRENADE MATERIALS:
RADIAL GRID ELEMENT OF HE
                                                   CM
TMD OF HE
                                                   G/CC
SPECIFIC HEAT OF HE
                                                   CAL/G/DEG K
THERMAL CONDUCTIVITY OF HE
                                                   CAL/CM/S/DEG K
THERMAL DIFFUSIVITY OF HE
                                                   SQ CM/S
RADIAL GRID ELEMENT OF STEEL
                                                   CM
TMD OF STEEL
                                                   G/CC
SPECIFIC HEAT OF STEEL
                                                   CAL/G/DEG K
THERMAL CONDUCTIVITY OF STEEL
                                                   CAL/CM/S/DEG K
THERMAL DIFFUSIVITY OF STEEL
                                                   SQ CM/S
  242
            PRINT 9 LINES WITH RHOHEO, PHE, PINT/1000000, XO, SIGKPSI, HSTRESS,
  243
           HSTRAIN, HSE, HSE*CALPJ, EE. RDX, EE. RDX*CALPJ, PWORK, PWORK*CALPJ
  244
           THUS
INITIAL CONDITIONS AT START OF FINAL CONSOLIDATION:
HE DENSITY
                                    G/CC
PRESSURE ON HE
                                    KPSI
                                                          MPA
HEIGHT OF HE COLUMN
                                    CM
MAX BODY HOOP STRESS
                                    KPSI
                                                          MPA
HOOP STRAIN IN BODY
HOOP STRAIN ENERGY
                                    JOULE
                                                          CAL
RDX ELASTIC ENERGY
                                    JOULE
                                                          CAL
CUM WORK BY PUNCH
                                    JOULE
                                                          CAL
  254
            SKIP 2 LINES
  255
           PRINT 7 LINES THUS
SIMULATED CONDITIONS DURING HE CONSOLIDATION
PUNCH
        HE
                 HE
                         MAX
                                  HOOP
                                           H STRN
                                                    PUNCH
                                                             THERM
                                                                     TEMP
                                                                              CYCLE
DISPL
                         STRESS
        DENS
                 PRESS
                                  STRAIN
                                          ENERGY
                                                    WORK
                                                            WORK
                                                                     RISE
                                                                              TIME
(CM)
       (G/CC)
                 (KPSI)
                          (KPSI)
                                  (MU/U)
                                           (CAL)
                                                    (CAL)
                                                             (CAL)
                                                                     (D K)
                                                                              (MS)
           WHILE XP < XPMAX, DO
  263
  264
  265
       ''INCREMENT DISPLACEMENT OF THE PUNCH.
       1 1
  266
  267
                ADD DELXP TO XP
```

```
268
             LET HOLDXP=XP
269
270
    ''CALC VOLUME AND DENSITY OF HE W/O ACCOUNTING FOR ADD'NL BODY STRAIN.
271
272
             LET RHOHE=MASSHE/AXHE/(XO-XP)
273
             LET PHE=PRESS(TMDHE, RHOHE)
274
             LET PINT_PA.KPSI*PHE
275
     ''CALCULATE INCREMENT OF WORK DONE BY PUNCH AT THIS PRESSURE.
276
277
278
             LET DELPWORK=0.000001 #AXHE PINT DELXP
             ADD DELPWORK TO PWORK
279
     11
280
281
     ''CALCULATE BODY STRAIN AT THIS PRESSURE.
282
283
             CALL STRS. STRAIN (PINT, POUT) YIELDING SIGPA, HSTRAIN
284
             LET SIGKPSI=SIGPA/PA.KPSI
285
286
     ''OBTAIN NEW CROSS SECTION, ACCOUNTING FOR CURRENT STRAIN.
287
288
             IFT AXHE=AXHEO*(1.0+HSTRAIN)**2
     . .
289
290
     "FIND THE ADDITIONAL PUNCH DISPLACEMENT WHICH PRESERVES HE DENSITY.
291
             LET XP=XO-MASSHE/AXHE/RHOHE
292
293
     ''ADD THE ADDITIONAL PUNCH WORK FOR THE ADDITIONAL DISPLACEMENT.
294
295
296
             DELPWORK=0.000001*AXHE*PINT*(XP-HOLDXP)
297
             ADD DELPWORK TO PWORK
     1 1
298
     ''UPDATE TIME.
299
300
             LET DELTAT=(XP-XPO) **RPSPEED(XP)
301
             LET DELT2=0.5*DELTAT
302
303
             ADD DELTAT TO TIME
304
     ''CALCULATE THE HOOP STRAIN ENERGY.
305
306
             LET HSE-STRN.ENER (PINT.POUT)
307
308
     ''CALCULATE THE ELASTIC STRAIN ENERGY IN THE RDX.
309
310
311
             J = EE.RDX=ESTRN.RDX(PINT)
312
     ''CALCULATE THE THERMAL ENERGY BALANCE.
313
314
315
             LET HENERGY=MAX.F(0.0, PWORK-HSE-EE.RDX-HTLOSS)
316
317
     ''CALCULATE THE THERMAL FLUX TO THE RDX.
318
319
             LET QDOT=CALPJ*(HENERGY-HENERGYO)/(MASS.RDX*SPS.RDX)/DELTAT
320
321
     ''CALCULATE AVG TEMP RISE IN HEATED HE.
322
323
             LET UHE=HENERGY*CALPJ/(HTCAP.RDX+HTCAP.STER)
     11
324
```

```
325
       ''CALCULATE HEAT LOSS TO WALL.
 326
               IF FLAGUHE NE 1
 327
 328
                   GO TO L3
               OTHERWISE ''GET BADIAL DIST OF TEMP IN HE AND IN STEEL WALL
 329
 330
               IF UCON.HE*DELTAT > 0.25
                   GO TO 1.3 ''TO AVOID NUMERICAL INSTABILITY
 331
 332
               OTHERWISE
               IF UCON.ST*DELTAT > 0.25
 333
                   GO TO L3 ''TO AVOID NUMERICAL INSTABILITY
 334
 335
               OTHERWISE
               LET TIMEU=TIME
 336
               LET OMEGA=QDOT*MASS.RDX*SPS.RDX/MASSHE ''HEAT GEN RATE/MASSHE
 337
 338
               FOR I=2 TO NHE-1 DO
                   LET RIA=DRHE#I
 339
 340
                   LET RI=RIA-DRHE
                   LET UHEV(I) = UHEOV(I) + DELTAT * UCON. HE/RI * (RIA * (UHEOV(I+1)
 341
 342
                   -UHEOV(I))-RI*(UHEOV(I)-UHEOV(I-1)))+DELTAT*OMEGA/SPHT.HE
                   LET UHEV(I)=MAX.F(0.0,UHEV(I))
 343
               LOOP ''OVER HE GRID
 344
               LET UHEV(1)=UHEV(2) ''AT CENTER OF HE
 345
               FOR I=2 TO NST-1 DO
 346
                   LET RIA=RINT+DRST*I
 347
                   LET RI=RIA-DRST
 348
                   LET USTV(I) = USTOV(I) + DELTAT*UCON.ST/RI*(RIA*(USTOV(I+1)
 349
                    -USTOV(I))-RI*(USTOV(I)-USTOV(I-1)))
 350
                   LET USTV(I)=MAX.F(0.0.USTV(I))
 351
               LOOP ''OVER STEEL GRID
 352
               LET USTV(NST)=USTV(NST-1) ''FOR INSULATED OUTER BOUNDARY
 353
  354
       ''AT HE-STEEL BOUNDARY, CONSERVE FLUX AND TEMPERATURE.
 355
  356
               LET UHEV(NHE) = (BETA#UHEV(NHE-1)+USTV(2)) #ROPB
  357
               LET USTV(1)=UHEV(NHE)
  358
  359
               LET DELHTLOSS=AWALL*TCOND.HE*(UHEV(NHE-1)-UHEV(NHE))/DRHE*
               DELTAT/CALPJ
  360
  361
       ''UPDATE INITIAL CONDITIONS.
  362
  363
               FOR I=1 TO NHE, LET UHEOV(I)=UHEV(I)
  364
  365
               FOR I=1 TO NST, LET USTOV(I)=USTV(I)
               GO TO L4
  366
               LET DELHTLOSS=AWALL#TCOND.HE#(UHE-USTEEL)/DELR.HE#DELTAT/CALPJ
  367
       'L3'
  368
       'L4'
               ADD DELHTLOSS TO HTLOSS
               LET USTEEL=HTLOSS/HTCAP.STEEL*CALPJ
  369
               ADD 1 TO K
  370
               IF FLAGUHE NE 1
  371
  372
                    GO TO L6
               OTHERWISE
  373
  374
               IF MOD.F(K, KUPRINI) NE O
  375
                   GO TO L6
  376
               OTHERWISE
               SKIP 1 LINE
  377
               PRINT 5 LINES WITH 1000*TIMEU
  378
               THUS
  379
TEMP DIST IN GRENADE AT TIME ***. *** MS
```

```
HE RADIAL
                                        TEMP INCR
                                                                            LOCATION
LOC (CM)
                                         (DEG K)
                                                                            INDEX
                                      FOR I=1 TO NHE DO
     385
     386
                                                LET J=I-1
     387
                                                LET R=DRHE*J
     388
                                                PRINT 1 LINE WITH R, UHEV(I), I
     389
                                                THUS
   **. ***
                                            *** ***
                                      LOOP ''OVER HE GRID
     391
                                      PRINT 4 LINES THUS
     392
WALL RADIAL
                                         TEMP INCR
                                                                            LOCATION
                                         (DEG K)
LOC (CM)
                                                                             INDEX
                                      FOR I=1 TO NST DO
     397
     398
                                                LET J=I-1
     399
                                                LET R=DRST#J+RINT
                                                 PRINT 1 LINE WITH R, USTV(I), J+NHE
     400
     401
    **.***
                                            *** ***
                                      LOOP ''OVER ST GRID
     403
                                      PRINT 2 LINES THUS
     404
                   'L6'
                                      IF FLAGU=0
     407
     408
                                                 GO TO L5
     409
                                       OTHERWISE ''GET TEMP PROFILE IN RDX PARTICLE
                                       IF UCON*DELT2 > 0.25
      410
      411
                                                 GO TO L5 ''TO AVOID NUMERICAL INSTABILITY
      412
                                       OTHERWISE
                                      LET U1V(2) = U0V(2) + DELT2 + UCON + (U0V(3) - U0V(2))
      413
      414
                                       LET U1V(1)=U1V(2) ''FOR ZERO COND AT CENTER OF PARTICLE
     415
                                       FOR I=3 TO NGRID-1 DO
      416
                                                 LET RI=(I-1) *DELTAR
      417
                                                 LET U1V(I) = U0V(I) + DELT2 = U1CON = ((U0V(I+1) + U0V(I-1))/RI + U1V(I) = U0V(I) + U1V(I) = U0V(I) + U1V(I) = U1V(I) 
      418
                                                  (UOV(I+1)-2.0*UOV(I)+UOV(I-1))/DELTAR)
                                       LOOP ''OVER GRID
      419
                                       LET U1V(NGRID)=U1V(NGRID-1)+U0CON*QDOT
      420
      421
      422
                   "MOVE NEXT HALF-TIME STEP.
      423
      424
                                       LET UV(2) = U1V(2) + DELT2 + UCON + (U1V(3) - U1V(2))
                                       LET UV(1)=UV(2) ''AT CENTER
      425
      426
                                       FOR I=3 TO NGRID-1 DO
      427
                                                  LET RI=(I-1)*DELTAR
      428
                                                 LET UV(I)=U1V(I)+DELT2#U1CON*((U1V(I+1)-U1V(I-1))/RI +
      429
                                                  (U1V(I+1)-2.0*U1V(I)+U1V(I-1))/DELTAR)
      430
                                       LOOP ''OVER GRID
      431
                                       LET UV(NGRID)=UV(NGRID-1)+U0CON*QDOT
      432
      433
                   ''UPDATE INITIAL TEMPERATURE DISTRIBUTION.
      434
                   . .
      435
                                       FOR I=1 TO NGRID, LET UOV(I)=UV(I)
      436
                   "PRINT TEMPERATURES AS REQUIRED.
      437
      438
```

```
439
               IF MOD.F(K, KUPRINT) NE O
  440
                   GO TO L5
  441
               OTHERWISE:
  442
               SKIP 1 LINE
  443
               PRINT 5 LINES WITH 1000*TIME
  444
               THUS
TEMP DISTRIBUTION IN RDX AT TIME *** MS
                TEMP INCR
RDX RADIAL
                               LOCATION
LOC (MICRON)
                (DEG K)
                               INDEX
               FOR I=1 TO NGRID DO
  450
  451
                   LET J=I-1
  452
                   LET R=10000.0*DELTAR*J
  453
                   PRINT 1 LINE WITH R,UV(I),J
  454
                   THUS
   ****
                 *** ***
  456
               LOOP ''OVER GRID
  457
               PRINT 2 LINES THUS
       'L5'
               IF PHE GE MAXKPSI
  460
                   GO TO L2
  461
  462
               OTHERWISE
  463
               IF MOD.F(K, KPRINT) NE O
  464
                   GO TO L1
  465
               OTHERWISE
               PRINT 1 LINE WITH XP, RHOHE, PHE, SIGKPSI, 10**6*HSTRAIN, CALPJ*
  466
  467
               HSE, CALPJ*PWORK, CALPJ*HENERGY, UHE, 1000*TIME
  468
 ****
  470
       'L1'
               LET XP0=XP
  471
               LET HENERGYO=HENERGY
  472
           LOOP ''OVER DISPLACEMENT
  473
               PRINT 1 LINE WITH XP, RHOHE, PHE, SIGKPSI, 10**6*HSTRAIN, CALPJ*
  474
               HSE, CALPJ*PWORK, CALPJ*HENERGY, UHE, 1000*TIME
  475
               THUS
  ***
                **,*** ***,** ****,* **,***
           PRINT 2 LINES THUS
  477
  480
       ''FOR THE RDX IN THE HE, GET THE STRAIN, STRAIN ENERGY, VOL RECOVERY,
  481
  482
       ''AND THE FINAL DENSITY OF THE HE.
  483
  484
           LET EPS.RDX=PINT/EY.RDX
  485
           LET EE.RDX=ESTRN.RDX(PINT) ''JOULE
  486
           LET DVOL. RDX=VOL. RDX#PINT/BM. RDX
  487
           LET DVOL.HE=VOL.RDX#EPS.RDX#(1.0+2.0*PR.RDX) ''UNI-AXIAL RELEASE
  488
           LET VOL.HE=MASSHE/RHOHE
  489
           LET RHOHE=RHOHE*VOL.HE/(VOL.HE+DVOL.HE) ''DENS AFTER RECOVERY
  490
  491
       ''HOOP STRESS ON THE OUTSIDE OF THE BODY AT AN INT PRESS OF PINT.
  492
  493
           LET A=0.01*RINT
  494
           LET B=0.01*ROUT
  495
           LET DENOM=B**2-A**2
```

```
496
           LET SIGOUT=(A**2*(PINT-POUT)+PINT*A**2-POUT*B**2)/DENOM
  497
           LET MAXLOAD=PI.C*(RINT/2.54) **2*PHE
  498
           PRINT 3 LINES WITH PINT/PA.KPSI, MAXLOAD, SIGOUT/PA.KPSI
  499
           THUS
                                    *** ***
MAX CONSOLIDATION PRESSURE
                                             KPSI
MAX CONSOLIDATION LOAD
                                             KLBF
                                    *** ***
HOOP STRESS ON BODY OUTSIDE
                                             KPSI
           PRINT 10 LINES WITH 10**6*EPS.RDX, VOL.RDX, DVOL.RDX, 100*DVOL.RDX/
  503
  504
           VOL. RDX, DVOL. HE, 100*DVOL. HE/VOL. HE, RHOHE, EE. RDX, CALPJ*EE. RDX,
  505
           HTLOSS, CALPJ*HTLOSS, USTEEL, 1000*TIME, QDOT
  506
           THUS
                         *****
STRAIN IN BULK RDX
                                   MU/U
VOLUME OF RDX SOLID
                                    CC
                                       (**. ***%)
RDX VOLUME INCREASE
                                    CC
                                        (** ***%)
1-AX HE VOL INCREASE
                                    CC
FINAL DENSITY OF HE
                                    G/CC
                                    JOULES
                                                 *** ***
RDX STRAIN ENERGY
                                    JOULES
CUM HEAT LOSS BY HE
                                                            CAL
TEMP RISE IN STEEL
                                   DEG K
THERMAL FLUX INTO RDX AT TIME ***. ** MS
                                                             CAL/SQ CM/S
  517
           SKIP 1 LINE
           IF FLAGUHE=1
  518
  519
                PRINT 5 LINES WITH 1000#TIMEU
  520
                THUS
TEMP DIST IN GRENADE AT TIME *** MS
                 TEMP INCR
HE RADIAL
                               LOCATION
LOC (CM)
                 (DEG K)
                               INDEX
  526
                FOR I=1 TO NHE DO
                    LET J=I-1
  527
  528
                    LET R=DRHE#J
  529
                    PRINT 1 LINE WITH R, UHEV(I), I
  530
                    THUS
 ** ***
                  ****
                LOOP ''OVER HE GRID
  532
                PRINT 4 LINES THUS
  533
WALL RADIAL
                 TEMP INCR
                               LOCATION
LOC (CM)
                 (DEG K)
                               INDEX
  538
                FOR I=1 TO NST DO
  539
                    LET J=I-1
  540
                    LET R=DRST#J+RINT
  541
                    PRINT 1 LINE WITH R, USTV(I), J+NHE
  542
                    THUS
  544
               LOOP ''OVER ST GRID
  545
                PRINT 2 LINES THUS
  548
           ALWAYS
  549
           STOP
  550
      END ''COMPACT
```

```
FUNCTION PRESS (TMD, RHO)
2
    ''FUNCTION CALCULATES THE TRANSIENT PRESSURE (KPSI) REQUIRED TO YIELD
    "HE DENSITY OF RHO (G/CC), WITH A THEORETICAL MAX HE DENSITY OF TMD.
    ''TRANSIENT PRESS IS THAT APPLIED DURING LOADING OP'NS OF THE M42/M46
    ''GRENADE. TMD OF THE HE IS ALSO GIVEN IN G/CC. THIS FUNCTION IS THE
    ''INVERSE OF RHOTRANS. CALCULATION OF THE PRESSURE FROM RHOTRANS IS
8
    ''OBTAINED ITERATIVELY VIA A NEWTON-RAPHSON METHOD.
 9
10
        LET ERR=0.0001 ''KPSI TOLERABLE ERROR
11
        LET PS= 0.8 RHO/(TMD-RHO) ''QUASI-STATIC PRESSURE FOR DENSITY RHO
        LET P1=1.52*PS
12
13
        LET P2=1.53*PS
14
        LET RHO1 = RHOTRANS(TMD, P1)
15
   'L1'LET RHO2=RHOTRANS(TMD, P2)
16
        IF ABS.F(P1-P2) LE ERR
17
            GO TO L3
18
        OTHERWISE
19
        LET P3=P2+(P2-P1)/(RHO2-RHO1)*(RHO-RHO2)
20
        LET P1=P2
21
        LET P2=P3
22
        LET RHO1=RHO2
23
        GO TO L1
24
    'L3'RETURN WITH P2
25
    END ''FUNCTION PRESS
    FUNCTION RHOTRANS (TMD, PRESS)
1
 2
    ''FUNCTION CALCULATES AVERAGE DENSITY OF THE HE WITHIN A GRENADE BODY
    ''FOR TRANSIENT LOADING CONDITIONS. ARGS: TMD IS THE THEORETICAL
 5
    "'Max DENSITY OF HE IN G/CC, AND PRESS IS THE PEAK PRESSURE IN KPSI.
    "PRESSURE GRADIENT WITHIN THE HE IS GIVEN BY THE COEF GRAD. PRS.
 7
8
        LET B=GRAD.PRS*PRESS
9
        LET ARG=(PRESS+0.8-B)/(PRESS+0.8)
10
        LET NORMRHO=1.0+0.8/B*LOG.E.F(ARG)
11
        RETURN WITH TMD*NORMRHO
12 END 'FUNCTION RHOTRANS
```

```
FUNCTION STRN. ENER (PINT, POUT)
   ''FUNCTION CALCULATES THE HOOP STRAIN ENERGY IN A CYLINDRICAL SLEEVE,
 3
   ''GIVEN INTERNAL PRESSURE PINT AND OUTSIDE PRESSURE POUT.
    ''HEIGHT, INTERNAL RADIUS, EXTERNAL RADIUS, AND YOUNG'S MODULUS FOR
6
    ''THE SLEEVE MATERIAL ARE TRANSMITTED AS GLOBAL VAHIABLES. INPUT
    ''ARGS ARE GIVEN IN PASCALS. HOOP-STRAIN ENERGY IS GIVEN IN JOULES.
    "'RESULT IS BASED ON THE LAME EQ'N FOR THICK-WALLED CYLINDERS.
9
10
        LET A=0.01*RINT
11
        LET B=0.01 #ROUT
12
        LET L=0.01*LBODY
        LET DENOM=B**2-A**2
13
        LET T1=A##2#B##2#(PINT-POUT)##2/2.0/DENOM
14
15
        LET T2=(PINT*A**2-POUT*B**2)**2/DENOM
16
        LET T3=2.0#A##2#B##2#(PINT-POUT)#(PINT#A##2-POUT#B##2)#LOG.E.F(B/A)
17
        /DENOM/DENOM
        RETURN WITH PI.C*L*(T1+T2+T3)/EYOUNG
18
19
   END ''STRN. ENER
    ROUTINE STRS. STRAIN GIVEN PINT, PCUT YIELDING SIGPA, HSTRAIN
 2
 3
    ''CALCULATES MAX HOOP STRESS AND HOOP STRAIN IN A CYLINDRICAL SLEEVE.
 Ц
    ''GIVEN INTERNAL PRESSURE PINT AND OUTSIDE PRESSURE POUT. THE SLEEVE
    "'INTERNAL RADIUS, EXTERNAL RADIUS, AND YOUNG'S MODULUS FOR SLEEVE
    ''MATERIAL ARE TRANSMITTED AS GLOBAL VARIABLES. THE LAME EQUATION FOR
 7
    ''THICK-WALLED CYLINDERS IS EVALUATED AT THE INT RADIUS.
                                                              DIMENSIONS
    ''ARE GIVEN IN CM. THEY ARE CONVERTED TO METERS. MAX STRESS (SIGPA)
 8
    ''IS GIVEN IN PASCALS. REF: TIMOSHENKO AND GOODIER, THEORY OF
    ''ELASTICITY, C. 1951.
10
    . .
11
12
        LET A=0.01 *RINT
        LET B=0.01 *ROUT
13
14
        LET DENOM=B*#2-A**2
        LET SIGPA=(PINT*(A**2+B**2)-2.0*B**2*POUT)/DENOM
15
16
        LET HSTRAIN=SIGPA/EYOUNG
17
        RETURN
18 END ''STRS.STRAIN
    FUNCTION ESTRN. RDX (PINT)
 1
 2
 3
    ''CALCULATES THE ELASTIC STRAIN ENERGY (JOULES) AS A FUNCTION OF PINT,
 4
    'THE PRESSURE (PA) ON THE HE . A LOCAL HIDROSTATIC STATE IS
    ''ASSUMED. VOLUME OF RDX CRYSTALS (VOL.RDX), AND VALUES OF YOUNG'S
    ''MODULUS (EY.RDX) AND BULK MODULUS (BM.RDX) FOR RDX ARE GLOBAL
 7
    ''INPUT ARGUMENTS. RDX VOLUME IS TREATED AS A CONSTANT.
    "PRESSURE GRADIENT IS ASSUMED WITHIN THE HE VOLUME.
                                                          PRESSURE VARIES
 9
    ''LINEARLY FROM A MAX OF PINT TO A MIN OF (PINT - GRAD.PRS*PINT).
10
   ''VALUE OF GRAD. PRS IS ASSIGNED IN THE MAIN PROGRAM AND TRANSMITTED
    ''GLOBALLY.
11
        LET EPS.RDX=PINT/EY.RDX
12
        LET COEF=1.0-GRAD.PRS+GRAD.PRS**2/3.0
13
14
        LET EE.RDX=0.0000005#VOL.RDX/BM.RDX#PINT##2#COEF
15
        RETURN WITH EE. RDX
    END ''FUNCTION ESTRN. RDX
```

```
FUNCTION RPSPEED (XP)
    ''CALCULATES THE RECRIPROCAL PUNCH SPEED IN S/CM AS A FUNCTION OF
3
    ''PUNCH DISPLACEMENT XP, IN CM. INITIAL, CONSTANT ACCELERATION OVER
4
    ''TRAVEL XCEL IS ASSUMED. A FINAL, CONSTANT DECELERATION IS ASSUMED
    ''OVER TRAVEL: XPMAX - DCEL TO XPMAX. SPEED IS CONSTANT WITHIN THE
6
    ''RANGE: XCEL TO XPMAX - DCEL. XPMAX IS A GLOBAL VARIABLE.
 7
8
9
        LET MAXSPEED=2.6 ''CM/S
        LET XCEL=0.02 ''CM
10
        LET DCEL=0.08 ''CM
11
12
        IF XP LE 0.0
13
           GO TO L1
14
        OTHERWISE
15
        IF FLAGU=1
16
            RETURN WITH 1.0/MAXSPEED
17
        OTHERWISE
18
        IF XP < XCEL
            LET SPEED=MAXSPEED#SQRT.F(XP/XCEL)
19
20
            RETURN WITH 1.0/SPEED
21
        OTHERWISE
22
        IF XP > XPMAX-DCEL
23
            LET SPEED=MAXSPEED*(1.0-SQRT.F((XP-XPMAX+DCEL)/DCEL))
24
            RETURN WITH 1.0/MAX.F(SPEED, 0.1)
25
        OTHERWISE
26
        RETURN WITH 1.0/MAXSPEED
    'L1'PRINT 1 LINE WITH XP THUS
ERROR. PUNCH DISPLACEMENT 1S OUT OF RANGE: XP = ....
29
        STOP
30 END ''FUNCTION RPSPEED
```

AMSMC-RDA-S

28 Aug 1985 Revised 24 Sep 85

MEMORANDUM FOR RECORD

SUBJECT: Predicted Bulk Densities of Comp A-5 and Comp A-4 As Functions of Peak Consolidation Pressure

1. Reference:

- a. Information Paper, SMCAR-ESM-M, 25 Ap 85, subject: Follow-up Questions on Comp A-5 Density.
- b. Ltr, SMCHO-QA, 2 July 85, subject: Density Measurements.
- c. Information Paper, SMCAR-ESM-M, 6 Aug 85, subject: Results from Comp A-5 Density Experiment.
- d. MFR, AMSMC-RDA-S, 9 Aug 85, subject: Some Observations About the Explosive Sensitivity of Comp A-5.
- e. MFR, DRSMC-SAS (R), 4 Nov 83, subject: Particle-Size Distribution of Nominal Class 1 RDX Before Incorporation and After Extraction From Extrusions of Comp C-4.

2. Background

In the Ref 1a paper SMCAR-ESM-M referenced some experimental results of consolidation of explosive Comp A-5 under quasi-static conditions. The applied pressure was permitted to remain at max value for several minutes in this quasi-static test, so that the explosive might be expected to exhibit hydrostatic behavior. The bulk density is given at several values of the peak consolidation pressure. The source of these data is indicated as George Ziegler, SMCAR-LCE. At the time these experiments were run the RDX granulation used in Comp A-5 was much coarser than it currently is. Consequently, there was some concern that the relationship between peak pressure and bulk density might be somewhat different now, using nominal Class 1 RDX, than it was then. Bill Fortune (SMCAR-ESM-M) directed Holston AAP to perform another set of quasi-static experiments on Comp A-5, having the current (nominal Class 1) granulation of RDX and, incidentally, reduced residual solvent level. I suggested a set of pressures at which these experiments were to be run. The range of pressure--7 to 60 kpsi--was greater than the range of Ziegler's data, with some matching pressure values. The results of the tests at Holston were transmitted in Ref These results are in good agreement with Ziegler's, as shown in Table 1. The average intra-batch standard deviation in the HSAAP data is 0.005 g/ml. An analytic model of the quasi-static, pressuredensity relationship is also shown, for comparison.

TABLE 1

COMPARISON OF COMP A-5 BULK DENSITY VERSUS PEAK QUASI-STATIC CONSOLIDATION PRESSURE

Entries are average density in g/ml.

	Data S	ource	
Pressure (kpsi)	ARDC Ziegler	Holston AAP 7/2/85 *	Calculated
7	1.61	1.599	1.597
10	1.65	1.633	1.648
15	1.69	1.686	1.690
17	1.70	1.692	1.700
24	1.72	1.725	1.722

- * Average of 3 batches of 5 samples per batch.
- ** density/TMD = p/(p + 0.8), where the theoretical max density (TMD) is 1.78 g/ml.

3. Effect of RDX Particle Size

The tabulated results support the assertion that the initial size of the RDX grains in Comp A-5 (within limits) does not have a measurable effect on the bulk density, obtained under quasistatic conditions. Ref ic also compares these experiments, but draws a stronger conclusion from them than the conclusion which I have made. In pgf 4 of Ref 1c one finds: "The conclusion then is that there appears to be no physical change to the explosive from the process changes." In view of the fact that bulk density is the only (!) physical quantity being measured here, it may be premature to make this assertion. In fact, evidence cited in Ref 1d indicates that grain size does affect explosive sensitivity in pressed, granular explosives. Further, at the range of density to which Comp A-5 is pressed in the M42/M46 grenade, there is indirect evidence (from tests on similar explosives) that grain fracture occurs during consolidation of these grenades. There is evidence (Ref 1e) of RDX grain fracture when Comp C-4 is extruded at pressures much lower than the 25 kpsi used on Comp A-5. Whether the change in explosive sensitivity due to different initial grain size is important is another issue. The fact is that the results in Table 1 do not address that issue.

4. Average Bulk Density of Comp A-5 After Transient Pressurization

As indicated in Ref 1c, the average bulk density of Comp A-5 under the transient conditions of loading in the M42/M46 grenade is less than the quasi-static value, given the same peak pressure. This fact is conjectured to result from a non-hydrostatic behavior of the explosive during the brief (120 millisec) time the pressure is at or near peak value. A gradient in pressure within the explosive would cause a corresponding gradient in density, yielding a smaller mean value. The compressibility of the explosive (locally) is given to a very good approximation by the quasistatic result:

$$rho/TMD = p/(p + 0.8),$$
 (1)

where the bulk density, rho, and TMD are given in, say, g/ml and where the peak consolidation pressure, p, is given in kpsi. For Comp A-5 the TMD is, practically, 1.78 g/ml. Now suppose that the max pressure reached in the explosive was not the same everywhere within the explosive volume. Specifically, as an approximation, assume a constant gradient of pressure with respect to volume, plotted from the max value on the punch face, where the hydrostatic value, pmax, should occur. Thus, the local pressure-volume relationship would have the form:

$$p = pmax - b x, (2)$$

with

b = pmax - pmin,

and where x is the non-dimensional volume coordinate, which varies from 0 to 1 over the volume.

To find the average bulk density (over the grenade internal volume), one must integrate the local density—a function of local pressure—over this volume. This corresponds to substituting equation (2) into (1), and integrating over the non-dimensional volume coordinate x. Formally, let the volume—average bulk density be denoted by E(rho). Then,

$$E(\text{rho})/\text{TMD} = x-\text{integral over } (0,1) \text{ of } :$$

$$(pmax - b x)/(pmax + 0.8 - b x) dx. \tag{3}$$

The result of this operation is a function with one arbitrary constant, b:

$$E(rho)/TMD = 1 + (0.8/b)ln((pmax+0.8-b)/(pmax+0.8)).$$
 (4)

In the limit when b approaches zero, equation (1) is produced, as expected. By examining the (Ref 1c) KSAAP experiments, it is found that a good fit to all results obtains when

$$b = 0.8 \text{ pmax}, \tag{5}$$

for values of pmax from 18 to 29 kpsi. With the above value of b substituted in equation (4), one has the following equation for average bulk density versus peak consolidation pressure.

$$E(rho)/TMD = 1 + (1/pmax)ln((0.2pmax+0.8)/(pmax+0.8))$$
 . (6)

5. Comparison of Experimental and Calculated Avg Bulk Density

Experimental (transient) average bulk density values, obtained in the KSAAP M42/M46 transient leading tests, were read from a graph provided in Ref 1c. A limited number of transient data points from MAAAP is also shown in Ref 1c. These data are compared (Table 2) with the calculated values of equation (6). The fit over the range appears to be within experimental error. Further, no trend in the residuals is apparent. On this basis one can assert that Comp A-5 falls far short of being an ideal fluid

during transient consolidation. A portion of the explosive volume may experience an effective pressure which is only 20 % of pmax in terms of its compressible response.

TABLE 2

COMPARISON OF EXPERIMENTAL AND THEORETICAL AVERAGE BULK DENSITIES OF COMP A-5 AFTER M42/M46 GRENADE LOADING

Peak Pressure		Average De	nsity (g/ml)	
(kpsi)	Exp (KSAAP)	Exp (MAAAP)	Calcul	ated
	VE 81/82	July 1985	Transient	Quasi-static
8		1.500	1.491	1.618
12		1.550	1.574	1.669
14		1.600	1.600	1.684
18	1.640		1.636	1.704
19	1.645		1.643	1.708
20	1.642		1.649	1.712
21	1.657		1.655	1.715
22	1.666		1.660	1.718
23	1.667	1.650	1.665	1.720
24	1.675		1.670	1.722
25	1.676		1.674	1.725
26	1.679		1.678	1.727
27	1.682		1.681	1.729
28	1.685		1.684	1.730
29	1.694		1.687	1.732
30			1.690	1.734

^{*} The quasi-static value, 1.78p/(p + 0.8), is offered for comparison.

6. Quasi-static Bulk Density of Comp A-4

It is noted that equation (1), which is used to calculate the quasi-static density of Comp A-5, is written in a non-dimensional form. Both sides of this equation are ratios of dimensional quantities. The only dimensional parameter which relates to compressibility of the explosive is the constant 0.8 kpsi. If this constant is essentially the same for Comp A-5 and Comp A-4, the quasi-static bulk density of Comp A-4 can be calculated from (1) using the TMD = 1.76 g/ml. This density is based on a formulation of Comp A-4 that has an addition of 3.0 % wax to RDX. The assumed density of wax is 0.91 to 0.92 g/ml. Since the formulation of Comp A-4 may vary slightly in percent wax, one should recalculate the TMD if the actual formulation is much different from that assumed. Variation in the TMD of A-4 is also contributed by different waxes. The bulk density values for Comp A-4 shown in Table 3. Comparable values of Comp A-5 are shown for comparison.

TABLE 3

COMPARISON OF BULK DENSITIES OF COMP A-5 AND COMP A-4

VERSUS PEAK QUASI-STATIC CONSOLIDATION PRESSURE

Entries are calculated bulk densities in g/ml.

Pressure	Explosive	Composition		
(kpsi)	A-5	A-4 with TMD	of:	
		1.760	1.770	
7	1.597	1.578	1.588	
8	1.618	1.599	1.609	
9	1.635	1.616	1.626	
10	1.648	1.628	1.539	
15	1.690	1.670	1.680	
17	1.700	1.680	1.590	
20	1.712	1.692	1.702	
24	1.722	1.702	1.713	
40	1.745	1.724	1.735	
60	1.757	1.736	1.747	

7. Conclusions and Recommendations

Mathematical models for the bulk density of Comp A-5 and Comp A-4 have been presented here. Under quasi-static pressurization the equation which relates bulk density to peak pressure has a very simple form. However, this result is shown to match experimental values of bulk density for Comp A-5 over the range of pressure from 7 kpsi to 60 kpsi. When the loading condition on the explosive is rather transient, with near-peak load persisting for only 120 milliseconds, the explosive does not have time to react in a hydrostatic manner. It has been shown that the average bulk densit; of the explosive in M42/M46 grenades is significantly lower than the quasi-static value obtained for a given pressure. A formula for the average bulk density, after transient loading, is derived here. This fo rula fits experimental results over the entire experimental pressure range for Comp A-5. A similar formula may be derived for Comp A-4, given some information about the nature of the pressure gradient which exists within the explosive during pressurization. Of course, the same assumption could be made for Comp A-4 as was made for Comp A-5. I have been reluctant to do this, since there is a large disparity in the portion binder in these two explosives. Comp A-5 has about half as much binder as Comp A-4. In the quasi-static case, the compressible behavior of these explosives is predicted to be much the same, relative to their respective TMDs. In the interest of expanding engineering knowledge, it is recommended (a) the bulk density predictions of Table 3 for Comp A-4 be verified empirically (as was done for Comp A-5), after quasi-static loading, and (b) the average density of explosive in M42 grenades be measured after transient loading using Comp A-4. Resources permitting, it would also be desirable to measure the local bulk density within the HE in M42 grenades at several locations. These data would support the assumption that the

TABLE 3

COMPARISON OF BULK DENSITIES OF COMP A-5 AND COMP A-4 VERSUS PEAK QUASI-STATIC CONSOLIDATION PRESSURE

Entries are calculated bulk densities in g/ml.

Pressure	Pressure Explosive Composition		
(kpsi)	A-5	A-4 with TMD	of:
		1.760	1.770
7	1.597	1.578	1.588
8	1.618	1.599	1.609
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15	1.690	1.670	1.680
17	1.700	1.680	1.690
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24	1.722	1.702	1.713
40	1.745	1.724	1.735
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Mathematical models for the bulk density of Comp A-5 and Comp A-4 have been presented here. Under quasi-static pressurization the equation which relates bulk density to peak pressure has a very simple form. However, this result is shown to match experimental values of bulk density for Comp A-5 over the range of pressure from 7 kpsi to 60 kpsi. When the loading condition on the explosive is rather transient, with near-peak load persisting for only 120 milliseconds, the explosive does not have time to react in a hydrostatic manner. It has been shown that the average bulk density of the explosive in M42/M46 grenades is significantly lower than the quasi-static value obtained for a given pressure. A formula for the average bulk density, after transient loading, is derived here. This formula fits experimental results over the entire experimental pressure range for Comp A-5. A similar formula may be derived for Comp A-4, given some information about the nature of the pressure gradient which exists within the explosive during pressurization. Of course, the same assumption could be made for Comp A-4 as was made for Comp A-5. I have been reluctant to do this, since there is a large disparity in the portion binder in these two explosives. Comp A-5 has about half as much binder as Comp A-4. In the quasi-static case, the compressible behavior of these explosives is predicted to be much the same, relative to their respective TMDs. In the interest of expanding engineering knowledge, it is recommended (a) the bulk density predictions of Table 3 for Comp A-4 be verified empirically (as was done for Comp A-5), after quasi-static loading, and (b) the average density of explosive in M42 grenades be measured after transient loading using Comp A-4. Resources permitting, it would also be desirable to measure the local bulk density within the HE in M42 grenades at several locations. These data would support the assumption that the

effective pressure exhibits a constant gradient within the HE volume. Local densities should be measured at three locations on the grenade axis, from the cone to the inside of the cap (dome). Also, at a midaxial station, three local radial positions should be sampled for bulk density at different polar angles.

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burge Soblember

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